

Jay Dry Cleaners

BRONX, NEW YORK

**Interim Remedial Measure Construction
Completion Report**

NYSDEC Site Number: C203119

Prepared for:

Mr. Suhak Lee

3211 Riverdale Avenue

Bronx, NY 10463

Prepared by:

Jason B Stewart Engineering, P.C.

228 Park Avenue South PMB 34864

New York, NY 10003

516-441-5800

APRIL 2026

CERTIFICATIONS

I, Jason Stewart, am currently a registered professional engineer licensed by the State of New York, I had primary direct responsibility for implementation of the remedial program activities, and I certify that the Interim Remedial Measure Workplan was implemented and that all construction activities were completed in substantial conformance with the Department-approved Interim Remedial Measure Workplan.

I certify that all documents generated in support of this report have been submitted in accordance with the DER's electronic submission protocols and have been accepted by the Department.

I certify that all data generated in support of this report have been submitted in accordance with the Department's electronic data deliverable and have been accepted by the Department.

I certify that all information and statements in this certification form are true. I understand that a false statement made herein is punishable as a Class "A" misdemeanor, pursuant to Section 210.45 of the Penal Law. I, Jason Stewart, of Jason B Stewart Engineering, P.C. , am certifying as Owner's Designated Site Representative

105632

NYS Professional Engineer #

April 3, 2026

Date

Jason Stewart

Signature



TABLE OF CONTENTS

| | |
|---|-----|
| CERTIFICATIONS..... | II |
| TABLE OF CONTENTS | III |
| LIST OF ACRONYMS | IV |
| CONSTRUCTION COMPLETION REPORT | 1 |
| 1.0 BACKGROUND | 1 |
| 2.0 DESCRIPTION OF REMEDIAL ACTIONS PERFORMED | 3 |
| 2.1 GOVERNING DOCUMENTS | 4 |
| 2.2 REMEDIAL PROGRAM ELEMENTS | 5 |
| 2.3 CONTAMINATED MATERIALS REMOVAL..... | 7 |
| 2.4 REMEDIAL PERFORMANCE/DOCUMENTATION SAMPLING | 7 |
| 2.6 DRAINAGE SYSTEM | 10 |
| 2.8 DEVIATIONS FROM THE INTERIM REMEDIAL WORK PLAN..... | 11 |
| LIST OF TABLES | 12 |
| LIST OF FIGURES | 13 |
| LIST OF APPENDICES | 14 |

LIST OF ACRONYMS

| Acronym | Definition |
|-------------------|---|
| BCA | Brownfield Cleanup Agreement |
| BCP | Brownfield Cleanup Program |
| CAMP | Community Air Monitoring Plan |
| CCR | Construction Completion Report |
| CIS-1,2-DCE | Cis-1,2-Dichloroethene |
| EC | Engineering Control |
| GAC | Granular Activated Carbon |
| HASP | Health and Safety Plan |
| IC | Institutional Control |
| IRM | Interim Remedial Measure |
| NYSDEC | New York State Department of Environmental Conservation |
| NYSDOH | New York State Department of Health |
| PCE | Tetrachloroethene |
| RAO | Remedial Action Objective |
| SC | Site Characterization |
| SCOS | Soil Cleanup Objectives |
| SMP | Site Management Plan |
| SSDS | Sub-Slab Depressurization System |
| SVE | Soil Vapor Extraction |
| TCE | Trichloroethene |
| ug/l | Microgram per liter |
| ug/m ³ | Microgram per cubic meter |
| VOCS | Volatile Organic Compounds |
| W.C. | Water column |

CONSTRUCTION COMPLETION REPORT

1.0 BACKGROUND

Mr. Suhak Lee entered into a Brownfield Cleanup Agreement (BCA) with the New York State Department of Environmental Conservation (NYSDEC) in July, 2019, to investigate and remediate a 0.26-acre property located in Wakefield section of Bronx, New York

1.1 Site Description

The Site is located at 840 East 233rd Street in the Wakefield section of Bronx, New York and identified as Block 4857, Lot 49 on the New York City Tax Map. Figure 1 contains a Site location map.

The Site is approximately 11,350 square feet (0.26 acres) in area and is improved with a one-story commercial building approximately 4,700 square feet (0.11 acres) in area with a partial basement and an asphalt-paved parking lot. The building is currently occupied for commercial use by Great Day Cleaners, which utilizes Exxon DF-2000, a non-chlorinated biodegradable dry-cleaning fluid.

The general vicinity of the Site is composed predominantly of residential and commercial uses. The site is bounded by a twelve-story residential apartment building to the north, two- and three-story residential multi-family buildings to the south, a one-story commercial building to the east, and a one-story gasoline filling station to the west.

The IRM activities were performed at the adjacent property to the east at 864-868 E 233rd St. The adjacent property to the east is identified as Block 4857, Lot 59 on the New York City Tax Map and contains 5,530 square foot building with storage in the basement and offices and a church on the first floor.

1.2 Previous Investigations

Previous environmental assessments performed between the years 2016 and 2021 at the subject property included soil vapor and indoor air analyses which identified a soil vapor encroachment condition and groundwater sampling which identified an area of chlorinated solvent impact along the eastern property boundary. Based on these findings, NYSDEC required that the Participant perform a soil vapor intrusion study at the neighboring building to the east (866 233rd Street). On January 22, 2021, ACT mobilized to 866 233rd Street to perform a soil vapor intrusion study consisting of one indoor ambient air sample and two sub-slab soil vapor samples. Analytical results for the ambient indoor air sample identified a concentration of trichlorethylene (TCE) at 7.2 ug/m³ and soil vapor concentrations of TCE of 26 ug/m³ and 28,000 ug/m³. Additionally, several chlorinated solvent compounds including tetrachloroethene (PCE), vinyl chloride, and cis-1,2-dichloroethene (DCE) were detected at elevated concentrations in soil vapor beneath this neighboring building. The results of the soil vapor intrusion study were evaluated using New York State Department of Health (NYSDOH) guidance and it was determined that mitigation measures were warranted to address the vapor intrusion at 866 233rd Street. Previous site assessments and investigations have also identified that the depth to the groundwater table beneath the basement of the building is approximately two to three feet.

2.0 DESCRIPTION OF REMEDIAL ACTIONS PERFORMED

Remedial activities completed at the Site were conducted in accordance with the NYSDEC-approved Interim Remedial Measure Workplan (IRMWP) for the Jay Dry Cleaners site (February, 2022) and IRMWP Addendum dated January 25, 2023.

On December 15, 2021, PWGC mobilized to the property along with ACT and NYSDEC to assess options for vapor intrusion mitigation measures at 866 233rd Street. During the site inspection, multiple locations were observed on the basement floor that were identified as potential pathways for soil vapor intrusion including three sump pits which were covered with sheet metal, trenches which appeared to be approximately two to three inches wide, and multiple gouged areas with exposed subsurface that ranged from one-inch to 6-inches in diameter. These trenches and sumps were installed by the current property owner to manage flooding issues. These identified potential pathways were sealed and covered under a March 2022 IRM Work Plan in August and September of 2022 when the trenches were channelized with fiberglass bottoms and covers were placed over the sump pits.

Based on the observations documented the site inspection and the shallow depth to the water table beneath the basement floor that prevents the installation of a subgrade mitigation system, such as a sub slab depressurization system (SSDS), PWGC proposed the following measures:

- Sealing of exposed soil within trenches and gouges with cement/grout.
- Installing permanent manhole covers above the sump pits.

These measures were proposed to act as a barrier against impacted vapors from encroaching into the indoor air of the basement at this building. Between August 22nd, 2022 and September 8th, 2022, the existing exposed dirt trenches were expanded and NDS Pro Series Channel Drains were installed in all slab penetrations throughout the basements of 864 and 866 East 233rd Street. Existing exposed brick sump pits were replaced with NDS 18" HDPE sealed catch basins to store water that is then discharged

through existing pumps and connections to the NYC sewer. At locations where trenches and gouges were observed, concrete and/or grout was applied to seal off the exposed subsurface where vapor pathways were able to infiltrate. At a minimum, the applied concrete or grout matched the thickness of approximately 6 inches of the existing slab at each location.

During the performance of the IRM, drainage of overland onsite stormwater was also determined to be required in addition to the sealing of the existing trenches. Therefore, a new stormwater system consisting of composite drains, subsurface piping, and sealed sumps was installed and provides a sealed cover system that minimizes exposure to contaminated soil and soil vapor beneath the site, while also providing a pathway for overland stormwater to exit the property.

The locations of the sealed drainage system are shown in Figure 1. A photo log is attached as Appendix B. Community air monitoring plan data sheets and field notes are attached in appendix D.

All deviations from the IRMWP are noted section 2.8 below.

2.1 GOVERNING DOCUMENTS

2.1.1 Site Specific Health & Safety Plan (HASP)

All remedial work performed under the IRMWP were in full compliance with governmental requirements, including Site and worker safety requirements mandated by Federal OSHA.

A site-specific Health and Safety Plan (HASP) was complied with for all remedial and invasive work performed at the Site.

2.1.2 Quality Assurance Project Plan (QAPP)

The QAPP was included as Appendix D of the Remedial Investigation Work Plan approved by the NYSDEC. The QAPP describes the specific policies, objectives, organization, functional activities, and quality assurance/quality control activities designed to achieve the project data quality objectives. During IRM performance and documentation sampling, laboratory analyses of groundwater, soil vapor, and indoor air

samples were conducted by Phoenix Environmental Laboratories, an NYSDOH, ELAP-certified laboratory. Laboratory analyses were conducted in accordance with USEPA methods and NYSDEC Analytical Services Protocol (ASP) B deliverable format. All data was provided in the Department's Electronic Data Deliverable (EDD) EQuIS format.

QA/QC procedures required by the NYSDEC ASP and USEPA methods were followed, including initial and continuing instrument calibrations and analysis of other samples (blanks, laboratory control samples, and matrix spikes/matrix spike duplicates). The laboratory provided sample bottles were pre-cleaned and preserved in accordance with the USEPA methods.

Data Usability Summary Reports (DUSRs) were prepared in accordance with DER-10 by a third-party validator, Jeri Rossi, CEAC. The DUSRs are included in Appendix G.

2.1.3 Community Air Monitoring Plan (CAMP)

Community air monitoring was performed during all invasive remedial activities, including installation drains. Monitoring consisted of trained personnel utilizing a hand-held Photoionization Detector (PID) and Personal Aerosol Dust Monitor (PDR) to monitor for the presence of organic vapors and Particulate matter.

All invasive remedial activities where the composite cover slab inside the building was disturbed where the basement was only occupied by OSHA-certified personnel.

Copies of all field data sheets relating to the CAMP are provided in electronic format in Appendix D.

2.2 REMEDIAL PROGRAM ELEMENTS

2.2.1 Contractors and Consultants

- Advanced Cleanup Technologies, Inc. installed the drainage system and was responsible for remedial performance/documentation sampling

- Action Remediation, Inc. was responsible for coordinating and transporting the soil waste generated
- PWGC was responsible for directing and supervising all work until 2023, when Jason B Stewart, PE of Jason B Stewart, Engineering P.C. became the certifying Engineer of Record.

2.2.2 Site Preparation

On December 15, 2021, PWGC mobilized to the property along with ACT and NYSDEC to assess options for mitigation measures at 866 East 233rd Street. During the site inspection, multiple locations were observed on the basement floor that were identified as potential pathways for soil vapor intrusion including three sump pits which were covered with sheet metal, trenches which appeared to be approximately two to three inches wide, and multiple gouged areas with exposed subsurface that ranged from one-inch to 6-inches in diameter.

Documentation of agency approvals required by the IRMWP is included in Appendix A.

2.2.3 General Site Controls

- The basement level of the 866 East 233rd Steet is secured by locking doors accessible only by tenants and owner.
- All waste derived from remedial activities were containerized in 55-gallon DOT-approved drums, stored onsite and transported offsite for disposal as manifested waste. Copies of the fully executed waste manifests are contained in Appendix C.

2.2.4 Nuisance controls

- Dust generated during interior construction was mitigated by ventilation fans with HEPA filters. All work was performed in the basement, which was only occupied by OSHA-certified personnel.
- All work vehicles were parked in the drycleaner parking lot or along side streets. No vehicle routing issues, or other complaints were received.

2.2.5 CAMP results

The Community Air Monitoring Plan provided for the collection and analysis of air samples during remedial construction activities to ensure proper protections were employed to protect workers and the neighboring community. Monitoring was performed from August 22nd, 2022, to September 8th, 2022, in compliance with the Community Air Monitoring Plan in the approved IRMWP. Monitoring included one station in the basement of 866 East 233rd Street, which consisted of a photoionization detector (PID) and PDR dust monitor.

Time Weight Average(TWA) Levels above 100 $\mu\text{g}/\text{m}^3$ of particulates were detected on days of sump pit installation on September 7 and 8, 2022.

On September 8th, 2022, a dust reading spike of 179 $\mu\text{g}/\text{m}^3$ was noted around 2 pm following the installation of a sealed sump pit. Work ceased for the day after the spike to suppress particulate matter concentrations.

Copies of all field data sheets relating to the CAMP are provided in electronic format in Appendix D.

2.3 CONTAMINATED MATERIALS REMOVAL

On March 1, 2023 and June 11, 2024, concrete and soil derived from the installation of the drainage system was containerized in four (4) 55-gallon DOT-approved drums, stored onsite and transported offsite for disposal by Action Remediation, Inc. as non-hazardous manifested waste. Waste manifests are enclosed in Appendix C.

A figure of the location of areas where excavations were performed is shown in Figure 2.

2.4 REMEDIAL PERFORMANCE/DOCUMENTATION SAMPLING

On January 22nd, 2021, two sub-slab soil samples (SS-5 and SS-6) and one indoor air sample (IA-3) were collected from the basement of 866 East 233rd Street

adjacent to the eastern boundary of the Site. Samples were collected over an 8-hr period.

On July 11, 2022, A Dri-Eaz HEPA 700 Air Scrubber with an activated carbon filter was installed in the basement of 864 and 866 East 233rd Street on July 11, 2022.. The Dri-Eaz HEPA 700 Air Scrubber has a variable speed fan that can produce up to 700cfm of flow with up to 6 air exchanges per hour in 7000 cu. ft. spaces. Specifications for the air scrubber are included in Appendix F. Following air testing on August 2, 2022, which showed minimal improvement and some increases of CVOC concentrations in indoor air with the air scrubber, the air scrubber was turned off on July 25, 2023 and maintenance was discontinued. Additional indoor air mitigation measures are being proposed as part of the BCP Remedial Action Workplan and will be submitted to NYSDEC and NYSDOH for review.

On August 2, 2022, two additional indoor air samples (IA-4 and IA-7) were collected from the basement of the adjacent building to the east approximately 1 month after the installation of an air scrubber. Indoor air samples IA-3, IA-4, and IA-7 were resampled on August 10, 2023 more than a month after the implementation of the interim remedial measure of sealing the basement floor. Samples were collected over an 24-hr period.

In 2021, PCE was detected in the indoor air sample IA-3 at 19 mg/m³. PCE was also detected in sub-slab soil vapor samples SS-5 and SS-6 at 73,000 mg/m³ and 400 mg/m³, respectively. According to Matrix B of the NYSDOH Guidance, mitigation is recommended for the concentrations of PCE detected in sub-slab soil vapor and indoor air at the Site. In 2022, after the installation of an air scrubber, PCE was detected in the indoor air samples IA-4, and IA-7 at concentrations of 22.2 mg/m³, and 19 mg/m³, respectively. These concentrations were all below the NYSDOH Air Guideline Value for PCE of 30 mg/m³. Following floor sealing mitigation measures, the concentrations of PCE in the indoor air samples IA-3, IA-4, and IA-7 had reduced to 17.7 mg/m³, 17.7 mg/m³, and 13.8 mg/m³, respectively.

In 2021, TCE was detected in the indoor air sample IA-3 at 7.2 mg/m³. TCE was also detected in sub-slab soil vapor samples SS-5 and SS-6 at 28,000 mg/m³ and 26

mg/m³, respectively. According to Matrix A of the NYSDOH Guidance, mitigation is recommended for the concentrations of TCE detected in sub-slab soil vapor and indoor air at the Site. In 2022, after the installation of an air scrubber, TCE was detected in the indoor air samples A-4, and IA-7 at concentrations of 12.4 mg/m³, and 10.2 mg/m³, respectively. These concentrations were all above the NYSDOH Air Guideline Value for TCE of 2 mg/m³. Following floor sealing mitigation measures, the concentrations of TCE in the indoor air samples IA-3, IA-4, and IA-7 had reduced to 7.63 mg/m³, 7.14 mg/m³, and 5.48 mg/m³, respectively.

In 2021, CIS12-DCE was detected in the indoor air sample IA-3 at 35 mg/m³. CIS12-DCE was also detected in sub-slab soil vapor samples SS-5 and SS-6 at 79,000 mg/m³ and 65 mg/m³, respectively. According to Matrix A of the NYSDOH Guidance, mitigation is recommended for the concentrations of CIS12-DCE detected in sub-slab soil vapor and indoor air at the Site. In 2022, after the installation of an air scrubber, CIS12-DCE was detected in the indoor air samples IA-4, and IA-7 at concentrations of 68.2 mg/m³, and 63.4 mg/m³, respectively. Following floor sealing mitigation measures, the concentrations of CIS12-DCE in the indoor air samples IA-3, IA-4, and IA-7 had reduced to 31.4 mg/m³, 30.8 mg/m³, and 21.3 mg/m³, respectively.

In 2021, Vinyl Chloride was detected in the indoor air sample IA-3 at 1 mg/m³. Vinyl Chloride was also detected in sub-slab soil vapor samples SS-5 and SS-6 at 7,900 mg/m³ and 1.9 mg/m³, respectively. According to Matrix C of the NYSDOH Guidance, mitigation is recommended for the concentrations of Vinyl Chloride detected in sub-slab soil vapor and indoor air at the Site. In 2022, after the installation of an air scrubber, Vinyl Chloride was detected in the indoor air samples IA-4, and IA-7 at concentrations of 4.65 mg/m³, and 4.85 mg/m³, respectively. Following floor sealing mitigation measures, the concentrations of Vinyl Chloride in the indoor air samples 0.23 mg/m³, 0.25 mg/m³, and below detection limits, respectively.

In 2021, Carbon Tetrachloride was detected at low levels in the indoor air sample IA-3 at 0.52 mg/m³. Carbon Tetrachloride was also detected in sub-slab soil vapor sample SS-6 at 0.32 mg/m³. According to Matrix A of the NYSDOH Guidance, no further action is recommended for the concentrations of Carbon Tetrachloride detected in

sub-slab soil vapor and indoor air at the Site. In 2022, after the installation of an air scrubber, Carbon Tetrachloride was detected in the indoor air samples IA-4, and IA-7 at concentrations of 0.53 mg/m³, and 0.51 mg/m³, respectively. Following floor sealing mitigation measures, the concentrations of Carbon Tetrachloride in the indoor air samples IA-3, IA-4, and IA-7 had reduced to 0.49 mg/m³, 0.5 mg/m³, and 0.47 mg/m³, respectively.

In 2021, 1,1-Dichloroethene (11-DCE) was not detected at in the indoor air samples IA-3. 11-DCE was also detected in sub-slab soil vapor sample SS-5 at 670 mg/m³. According to Matrix A of the NYSDOH Guidance, mitigation is recommended for the concentrations of 11-DCE detected in sub-slab soil vapor at the Site. In 2022, after the installation of an air scrubber, 11-DCE was detected in the indoor air sample IA-4 at a concentration of 0.22 mg/m³ but was not detected in IA-7. Following floor sealing mitigation measures, 1,1-Dichloroethane was not detected in any of the indoor air samples.

A table and figure summarizing all air samples is included in Table 1 and Figure 3, respectively. Laboratory Reports are attached in Appendix E.

2.6 DRAINAGE SYSTEM

Exposure to remaining contamination in soil/fill at the site is prevented by a sealed drainage cover system placed over 864 and 866 East 233rd Street. The NDS Pro Series Channel Drains are made of Polyolefin plastic and provide a sealed pathway for overland water entering to the basement to be routed to catch basins installed in the basement floor. Existing exposed brick sump pits were replaced with NDS 18" HDPE sealed catch basins to store water that is then discharged through existing pumps and connections to the NYC sewer. The newly installed drains and catch basins were sealed to the existing concrete slab using non-shrink concrete grout. The resultant system is a completed capped drainage system that minimizes exposure to contaminated soil and soil vapor beneath the site.

The digital photo log is included in electronic format in Appendix B. Specifications for the sealed drains and sumps are included in Appendix F.

2.8 DEVIATIONS FROM THE INTERIM REMEDIAL WORK PLAN

During the performance of the IRM, drainage of overland onsite stormwater was also determined to be required in addition to the sealing of the existing trenches. Therefore, a new stormwater system consisting of composite drains, subsurface piping, and sealed sumps was installed and provides a sealed cover system that minimizes exposure to contaminated soil and soil vapor beneath the site, while also providing a pathway for overland stormwater to exit the property.

A Dri-Eaz HEPA 700 Air Scrubber with an activated carbon filter was installed in the basement of 864 and 866 East 233rd Street on July 11, 2022.. The Dri-Eaz HEPA 700 Air Scrubber has a variable speed fan that can produce up to 700cfm of flow with up to 6 air exchanges per hour in 7000 cu. ft. spaces. Specifications for the air scrubber are included in Appendix F. Following air testing on August 2, 2022, which showed minimal improvement and some increases of CVOC concentrations in indoor air with the air scrubber, the air scrubber was turned off on July 25, 2023 and maintenance was discontinued. Additional indoor air mitigation measures are being proposed as part of the BCP Remedial Action Workplan and will be submitted to NYSDEC and NYSDOH for review.

Following approval of the IRM addendum letter, dewatering wells DW-1 and DW-2 were installed in the basement of 866 East 233rd Street. The dewatering wells were sampled on March 9, 2023. DW-2 contained cis-1,2-Dichloroethylene at 22,100 ug/L and vinyl chloride at 4,640 ug/L. The results of this were included in the RIR and RAWP. To avoid continually dewatering, treating and drawing in the contaminated water, it was concluded the dewatering be better handled during the remedial action when treatment with ISCO injections is proposed.

LIST OF TABLES

1. Volatile Organic Compounds in Soil Vapor (ug/m³)
2. Volatile Organic Compounds in Indoor and Outdoor Air (ug/m³)
3. Volatile Organic Compounds in Groundwater (ug/L)
4. NYSDOH Air Matrices(Updated February 2024)

LIST OF FIGURES

1. Project Site Map
2. As-Built Site Diagram
3. Soil Vapor Intrusion Spider Diagram
4. Groundwater Spider Diagram

LIST OF APPENDICES

- A. NYSDEC Approvals of Substantive Technical Requirements
- B. Project Photo Log
- C. Waste Manifests
- D. CAMP Field Data Sheets and Air Monitoring Data
- E. Raw Analytical Laboratory Data
- F. Engineering Control Specifications
- G. Data Usability Summary Reports

Tables

Table 1
 Volatile Organic Compounds in Soil Vapor (ug/m3)
 EPA Method TO15

BCP Site No. C203119

| Sample ID | SV-1 | SV-2 | SV-3 | SS-1 | SS-2 | SS-3 | SS-4 | SS-5 | SS-6 | SS-12 | SS-15 | SS-18 | | | |
|---|-------------|------------|------------|------------|------------|------------|------------|------------|------------|---------|------------|---------|---|--------|----|
| York & Phoenix ID | 21D0660-01 | 21D0660-02 | 21D0660-03 | 20H0017-01 | 20H0017-02 | 20H0017-04 | 20H0017-05 | 21A0950-01 | 21A0950-02 | CM85136 | 22F1004-01 | CP07985 | | | |
| Sampling Date | 4/13/21 | 4/13/21 | 4/13/21 | 7/31/20 | 7/31/20 | 7/31/20 | 7/31/20 | 1/22/21 | 1/22/21 | 11/9/22 | 6/16/22 | 9/21/23 | | | |
| Client Matrix | Soil Vapor | | Soil Vapor | | Soil Vapor | | Soil Vapor | | Soil Vapor | | Soil Vapor | | | | |
| Compound | CAS Number | Result | Q | Result | Q | Result | Q | Result | Q | Result | RI | Result | Q | Result | RI |
| Volatile Organics, EPA TO15 Full List | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | |
| Dilution Factor | | 1.546 | | 1.753 | | 1.702 | | 45.37 | | 86.88 | | 662.14 | | 16.87 | |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | 1.100 | U | 1.200 | U | 1.200 | U | 31 | U | 12 | U | 12 | U | 16 | U |
| 1,1,1-Trichloroethane | 71-55-6 | 0.840 | U | 0.960 | U | 0.930 | U | 25 | U | 12 | U | 9.8 | U | 13 | U |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | 1.100 | U | 1.200 | U | 1.200 | U | 31 | U | 12 | U | 12 | U | 16 | U |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 76-13-1 | 1.200 | U | 1.300 | U | 1.300 | U | 35 | U | 17 | U | 13 | U | 18 | U |
| 1,1,2-Trichloroethane | 79-00-5 | 0.840 | U | 0.960 | U | 0.930 | U | 25 | U | 12 | U | 9.8 | U | 13 | U |
| 1,1-Dichloroethane | 75-34-3 | 0.630 | U | 0.710 | U | 0.690 | U | 18 | U | 8.8 | U | 6.8 | U | 7.3 | U |
| 1,1-Dichloroethylene | 75-35-4 | 0.150 | U | 0.170 | U | 0.170 | U | 4.5 | U | 2.2 | U | 1.7 | U | 1.8 | U |
| 1,2,4-Trichlorobenzene | 120-82-1 | 1.100 | U | 1.300 | U | 1.300 | U | 34 | U | 16 | U | 13 | U | 17 | U |
| 1,2,4-Trimethylbenzene | 95-63-6 | 3.800 | D | 5.300 | D | 3 | D | 22 | U | 11 | U | 8.3 | D | 8.9 | J |
| 1,2-Dibromoethane | 106-93-4 | 1.200 | U | 1.300 | U | 1.300 | U | 35 | U | 17 | U | 13 | U | 14 | U |
| 1,2-Dichlorobenzene | 95-50-1 | 0.930 | U | 1.100 | U | 1 | U | 27 | U | 13 | U | 10 | U | 11 | U |
| 1,2-Dichloroethane | 107-06-2 | 0.630 | U | 0.710 | U | 0.690 | U | 18 | U | 8.8 | U | 6.8 | U | 7.3 | U |
| 1,2-Dichloropropane | 78-87-5 | 0.710 | U | 0.810 | U | 0.790 | U | 21 | U | 10 | U | 7.8 | U | 8.3 | U |
| 1,2-Dichlorotetrafluoroethane | 76-14-2 | 1.100 | U | 1.200 | U | 1.200 | U | 32 | U | 15 | U | 12 | U | 13 | U |
| 1,3,5-Trimethylbenzene | 108-67-8 | 1.300 | D | 2 | D | 0.840 | J | 22 | J | 11 | U | 8.3 | U | 8.9 | U |
| 1,3-Butadiene | 106-99-0 | 1 | U | 1.400 | D | 1.100 | U | 30 | U | 14 | U | 11 | U | 12 | U |
| 1,3-Dichlorobenzene | 541-73-1 | 0.930 | U | 1.100 | U | 1 | U | 27 | U | 13 | U | 10 | U | 11 | U |
| 1,3-Dichloropropane | 142-28-9 | 0.710 | U | 0.810 | U | 0.790 | U | 21 | U | 10 | U | 7.8 | U | 8.3 | U |
| 1,4-Dichlorobenzene | 106-46-7 | 0.930 | U | 1.100 | U | 1 | U | 27 | U | 13 | U | 10 | U | 11 | U |
| 1,4-Dioxane | 123-91-1 | 1.100 | U | 1.300 | U | 1.200 | U | 33 | U | 16 | U | 12 | U | 13 | U |
| 2-Butanone | 78-93-3 | 1.300 | D | 20 | D | 1.400 | D | 13 | U | 6.4 | U | 25 | D | 11 | D |
| 2-Hexanone | 591-78-6 | 1.300 | J | 1.400 | J | 1.400 | U | 37 | U | 18 | U | 14 | U | 15 | U |
| 3-Chloropropene | 107-05-1 | 2.400 | U | 2.700 | U | 2.700 | U | 71 | U | 34 | U | 26 | U | 28 | U |
| 4-Methyl-2-pentanone | 108-10-1 | 0.630 | U | 3 | D | 0.700 | U | 19 | U | 8.9 | U | 6.9 | U | 7.4 | U |
| Acetone | 67-64-1 | 11 | D | 65 | D | 9.900 | D | 620 | D | 200 | D | 210 | D | 93 | D |
| Acrylonitrile | 107-13-1 | 0.340 | U | 0.380 | U | 0.370 | U | 9.8 | U | 4.7 | U | 3.7 | U | 3.9 | U |
| Benzene | 71-43-2 | 1.300 | D | 26 | D | 0.540 | U | 35 | D | 6.9 | U | 6.5 | D | 12 | D |
| Benzyl chloride | 100-44-7 | 0.800 | U | 0.910 | U | 0.880 | U | 23 | U | 11 | U | 8.7 | U | 9.3 | U |
| Bromodichloromethane | 75-27-4 | 1 | U | 1.200 | U | 1.100 | U | 30 | U | 15 | U | 11 | U | 12 | U |
| Bromofrom | 75-25-2 | 1.600 | U | 1.800 | U | 1.800 | U | 47 | U | 22 | U | 17 | U | 19 | U |
| Bromomethane | 74-83-9 | 0.600 | U | 0.680 | U | 0.660 | U | 18 | U | 8.4 | U | 6.6 | U | 7 | U |
| Carbon disulfide | 75-15-0 | 2.100 | D | 67 | D | 1.300 | D | 69 | D | 22 | D | 8.9 | D | 13 | D |
| Carbon tetrachloride | 56-23-5 | 0.490 | D | 0.280 | U | 0.640 | D | 7.1 | U | 3.4 | U | 2.7 | U | 2.8 | U |
| Chlorobenzene | 108-90-7 | 0.710 | U | 0.810 | U | 0.780 | U | 21 | U | 10 | U | 7.8 | U | 8.3 | U |
| Chloroethane | 75-00-3 | 0.410 | U | 0.460 | U | 0.450 | U | 12 | U | 5.7 | U | 4.5 | U | 4.8 | U |
| Chloroform | 67-66-3 | 8.200 | D | 4 | D | 67 | D | 22 | U | 23 | D | 8.3 | U | 8.8 | U |
| Chloromethane | 74-87-3 | 0.320 | U | 0.360 | U | 0.350 | U | 9.4 | U | 4.5 | U | 3.5 | U | 3.7 | U |
| cis-1,2-Dichloroethylene | 156-59-2 | 0.150 | U | 0.420 | D | 0.170 | U | 4.5 | U | 1.600 | D | 1.7 | U | 5.7 | D |
| cis-1,3-Dichloropropylene | 10061-01-5 | 0.700 | U | 0.800 | U | 0.770 | U | 21 | U | 9.9 | U | 7.7 | U | 8.2 | U |
| Cyclohexane | 110-82-7 | 0.530 | U | 3.100 | D | 0.590 | U | 7,400 | D | 51 | D | 5.8 | U | 6.2 | U |
| Dibromochloromethane | 124-48-1 | 1.300 | U | 1.500 | U | 1.400 | U | 39 | U | 19 | U | 14 | U | 15 | U |
| Dichlorodifluoromethane | 75-71-8 | 2.300 | D | 2.300 | D | 2.300 | D | 22 | U | 11 | U | 8.4 | U | 8.9 | U |
| Ethyl acetate | 141-78-6 | 1.100 | U | 1.300 | U | 1.200 | U | 33 | U | 16 | U | 12 | U | 13 | U |
| Ethyl Benzene | 100-41-4 | 10 | D | 3.300 | D | 0.810 | D | 61 | D | 9.4 | J | 7.3 | U | 7.8 | J |
| Hexachlorobutadiene | 87-68-3 | 1.600 | U | 1.900 | U | 1.800 | U | 48 | U | 23 | U | 18 | U | 19 | U |
| Isopropanol | 67-63-0 | 0.760 | J | 1.400 | D | 0.840 | J | 22 | U | 11 | U | 8.3 | U | 8.9 | U |
| Methyl Methacrylate | 80-62-6 | 0.630 | J | 0.720 | U | 0.700 | J | 19 | U | 8.9 | U | 6.9 | J | 7.4 | U |
| Methyl tert-butyl ether (MTBE) | 1634-04-4 | 0.560 | U | 0.630 | U | 0.610 | U | 16 | U | 7.8 | U | 6.1 | U | 6.5 | U |
| Methylene chloride | 75-09-2 | 1.100 | D | 1.900 | D | 1.200 | J | 32 | J | 15 | U | 12 | J | 13 | J |
| n-Heptane | 142-82-5 | 0.630 | J | 11 | D | 0.700 | U | 19 | U | 8.9 | U | 44 | D | 21 | D |
| n-Hexane | 110-54-3 | 0.930 | D | 11 | D | 0.780 | D | 1,300 | D | 42 | D | 30 | D | 15 | D |
| o-Xylene | 95-47-6 | 7 | D | 5.100 | D | 1.300 | D | 20 | U | 9.4 | J | 7.3 | J | 7.8 | J |
| p- & m- Xylenes | 179601-23-1 | 22 | D | 11 | D | 3.900 | D | 39 | U | 19 | J | 15 | U | 16 | U |
| p-Ethyltoluene | 622-96-8 | 3 | D | 5.100 | D | 2 | D | 27 | D | 11 | U | 8.3 | J | 8.9 | U |
| Propylene | 115-07-1 | 2.800 | D | 36 | D | 1.800 | D | 370 | D | 32 | D | 21 | D | 13 | D |
| Styrene | 100-42-5 | 0.660 | J | 0.750 | U | 0.730 | U | 19 | U | 9.3 | U | 7.2 | U | 7.7 | U |
| Tetrachloroethylene | 127-18-4 | 95 | D | 80 | D | 12 | D | 130 | D | 9,700 | D | 15,000 | D | 6,400 | D |
| Tetrahydrofuran | 109-99-9 | 1.100 | D | 1 | U | 1 | U | 27 | U | 13 | U | 10 | U | 14 | U |
| Toluene | 108-88-3 | 17 | D | 19 | D | 1.600 | D | 56 | D | 16 | D | 11 | D | 21 | D |
| trans-1,2-Dichloroethylene | 156-60-5 | 0.610 | U | 0.700 | U | 0.670 | U | 31 | D | 25 | D | 6.7 | U | 7.2 | U |
| trans-1,3-Dichloropropylene | 10061-02-6 | 0.700 | U | 0.800 | U | 0.770 | U | 21 | U | 9.9 | U | 7.7 | U | 8.2 | U |
| Trichloroethylene | 79-01-6 | 7.300 | D | 2.900 | D | 0.230 | U | 6.1 | U | 1,500 | D | 42 | D | 77 | D |
| Trichlorofluoromethane (Freon 11) | 75-69-4 | 1.600 | D | 1.300 | D | 1.300 | D | 25 | U | 12 | U | 9.5 | U | 10 | U |
| Vinyl acetate | 108-05-4 | 0.540 | U | 0.620 | U | 0.600 | U | 16 | U | 7.6 | U | 6 | U | 6.4 | U |
| Vinyl bromide | 593-60-2 | 0.680 | U | 0.770 | U | 0.740 | U | 20 | U | 9.5 | U | 7.4 | U | 7.9 | U |
| Vinyl Chloride | 75-01-4 | 0.200 | U | 0.220 | U | 0.220 | U | 59 | D | 2.8 | U | 2.2 | U | 2.3 | U |
| 4-Isopropyltoluene | 99-87-6 | NT | | NT | | NT | | NT | | NT | | NT | | NT | |
| Ethanol | 64-17-5 | NT | | NT | | NT | | NT | | NT | | NT | | NT | |
| Isopropylbenzene | 98-82-8 | NT | | NT | | NT | | NT | | NT | | NT | | NT | |
| n-Butylbenzene | 104-51-8 | NT | | NT | | NT | | NT | | NT | | NT | | NT | |
| sec-Butylbenzene | 135-98-8 | NT | | NT | | NT | | NT | | NT | | NT | | NT | |

NOTES:
 Yellow highlight = NYSDOH Indoor Air Quality Guideline exceedance
Q is the Qualifier Column with definitions as follows:
 D=result is from an analysis that required a dilution
 J=analyte detected at or above the MDL (method detection limit) but below the RL (Reporting Limit) - data is estimated
 U=analyte not detected at or above the level indicated
 B=analyte found in the analysis batch blank
 E=result is estimated and cannot be accurately reported due to levels encountered or interferences
 NT=this indicates the analyte was not a target for this sample
 ~this indicates that no regulatory limit has been established for this analyte
 RL = Reporting Limit

Table 2
Volatile Organic Compounds in Indoor and Outdoor Air (ug/m3)
 EPA Method TO15

BCP Site No. C203119

| Sample ID York & Phoenix ID | Sampling Date | NYSDOH Air Guideline Values | IA-1 20H0017-03 7/31/20 | | IA-2 20H0017-06 7/31/20 | | IA-3 21A0950-03 1/22/21 | | IA-4 CL97021 8/2/22 | | IA-7 CL97022 8/2/22 | | IA-3 21A0950-03 8/10/23 | | IA-4 CL97021 8/10/23 | | IA-7 CL97022 8/10/23 | | IA-11 CM85137 11/9/22 | | IA-14 22F1004-03 6/17/22 | | IA-17 CP07983 9/21/23 | | OA-1 20H0017-07 7/31/20 | | OA-2 CM85138 11/9/22 | | OA-3 22F1004-02 6/17/22 | | OA-4 CP07984 9/21/23 | | | | |
|---|---------------|--------------------------------|-------------------------------|--------------------|-------------------------------|--------------------|-------------------------------|--------------------|---------------------------|--------------------|---------------------------|--------------------|-------------------------------|--------------------|----------------------------|--------------------|----------------------------|--------------------|-----------------------------|--------------------|--------------------------------|--------------------|-----------------------------|--------------------|-------------------------------|--------------------|----------------------------|--------------------|-------------------------------|---------------------|----------------------------|---------------------|-------|-------------|------|
| | | | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Indoor Ambient Air | Outdoor Ambient Air | Outdoor Ambient Air | Outdoor Ambient Air | Outdoor Ambient Air | | | |
| Compound | CAS Number | | Result | Q | Result | Q | Result | Q | Result | RL | Result | RL | Result | RL | Result | RL | Result | RL | Result | RL | Result | RL | Result | RL | Result | RL | Result | RL | Result | RL | Result | RL | | | |
| Volatile Organics, EPA TO15 Full List | | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | ug/m3 | | | | |
| Dilution Factor | | | 0.975 | | 0.947 | | 1.037 | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | ~ | 0.67 | U | 0.65 | U | 0.71 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.64 | U | <1.00 | 1.00 | 0.8 | U | <1.00 | 1.00 | | | | | |
| 1,1,1-Trichloroethane | 71-55-6 | ~ | 0.53 | U | 0.52 | U | 0.57 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.51 | U | <1.00 | 1.00 | 0.64 | U | <1.00 | 1.00 | | | | | |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | ~ | 0.67 | U | 0.65 | U | 0.71 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.64 | U | <1.00 | 1.00 | 0.8 | U | <1.00 | 1.00 | | | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 76-13-1 | ~ | 0.97 | D | 0.73 | U | 0.79 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.84 | U | <1.00 | 1.00 | 0.72 | U | <1.00 | 1.00 | 0.89 | U | <1.00 | 1.00 | |
| 1,1,2-Trichloroethane | 79-00-5 | ~ | 0.53 | U | 0.52 | U | 0.57 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.6 | U | <1.00 | 1.00 | 0.51 | U | <1.00 | 1.00 | 0.64 | U | <1.00 | 1.00 | |
| 1,1-Dichloroethane | 75-34-3 | ~ | 0.39 | U | 0.38 | U | 0.42 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.44 | U | <1.00 | 1.00 | 0.38 | U | <1.00 | 1.00 | 0.47 | U | <1.00 | 1.00 | |
| 1,1-Dichloroethylene | 75-35-4 | ~ | 0.097 | U | 0.094 | U | 0.1 | U | 0.22 | 0.20 | <0.20 | 0.20 | <0.20 | 0.20 | <0.20 | 0.20 | <0.20 | 0.20 | <0.20 | 0.20 | <0.20 | 0.20 | 0.11 | U | <0.20 | 0.20 | 0.093 | U | <0.20 | 0.20 | 0.12 | U | <0.20 | 0.20 | |
| 1,2,4-Trichlorobenzene | 120-82-1 | ~ | 0.72 | U | 0.7 | U | 0.77 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.81 | U | <1.00 | 1.00 | 0.69 | U | <1.00 | 1.00 | 0.87 | U | <1.00 | 1.00 | |
| 1,2,4-Trimethylbenzene | 95-63-6 | ~ | 2.3 | D | 1.2 | D | 1 | D | 5.94 | 1.00 | 6.04 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 7.5 | D | <1.00 | 1.00 | 0.6 | D | <1.00 | 1.00 | 0.57 | J | <1.00 | 1.00 | |
| 1,2-Dibromoethane | 106-93-4 | ~ | 0.75 | U | 0.73 | U | 0.8 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.84 | U | <1.00 | 1.00 | 0.72 | U | <1.00 | 1.00 | 0.9 | U | <1.00 | 1.00 | |
| 1,2-Dichlorobenzene | 95-50-1 | ~ | 0.59 | U | 0.57 | U | 0.62 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.66 | U | <1.00 | 1.00 | 0.56 | U | <1.00 | 1.00 | 0.7 | U | <1.00 | 1.00 | |
| 1,2-Dichloroethane | 107-06-2 | ~ | 0.51 | D | 0.38 | U | 0.42 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.44 | U | <1.00 | 1.00 | 0.38 | U | <1.00 | 1.00 | 0.47 | U | <1.00 | 1.00 | |
| 1,2-Dichloropropane | 78-87-5 | ~ | 0.45 | U | 0.44 | U | 0.48 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.51 | U | <1.00 | 1.00 | 0.43 | U | <1.00 | 1.00 | 0.54 | U | <1.00 | 1.00 | |
| 1,2-Dichlorotetrafluoroethane | 76-14-2 | ~ | 0.68 | U | 0.66 | U | 0.72 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.77 | U | <1.00 | 1.00 | 0.65 | U | <1.00 | 1.00 | 0.82 | U | <1.00 | 1.00 | |
| 1,3,5-Trimethylbenzene | 108-67-8 | ~ | 0.77 | D | 0.47 | D | 0.51 | J | 1.98 | 1.00 | 1.99 | 1.00 | 2.42 | 1.00 | 3.04 | 1.00 | 6.04 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 2.3 | D | <1.00 | 1.00 | 0.46 | U | <1.00 | 1.00 | 0.57 | U | <1.00 | 1.00 | |
| 1,3-Butadiene | 106-99-0 | ~ | 0.65 | U | 0.63 | U | 0.69 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.73 | U | <1.00 | 1.00 | 0.62 | U | <1.00 | 1.00 | 0.77 | U | <1.00 | 1.00 | |
| 1,3-Dichlorobenzene | 541-73-1 | ~ | 0.59 | U | 0.57 | U | 0.62 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.66 | U | <1.00 | 1.00 | 0.56 | U | <1.00 | 1.00 | 0.7 | U | <1.00 | 1.00 | |
| 1,3-Dichloropropane | 142-28-9 | ~ | 0.45 | U | 0.44 | U | 0.48 | U | NT | | NT | | NT | | NT | | NT | | NT | | NT | | 0.51 | U | | 0.43 | U | NT | | 0.54 | U | NT | | | |
| 1,4-Dichlorobenzene | 106-46-7 | ~ | 1.2 | D | 0.57 | U | 0.62 | U | 1.06 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 1.2 | D | <1.00 | 1.00 | 0.56 | U | <1.00 | 1.00 | 0.7 | U | <1.00 | 1.00 | |
| 1,4-Dioxane | 123-91-1 | ~ | 0.7 | U | 0.68 | U | 0.75 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.79 | U | <1.00 | 1.00 | 0.67 | U | <1.00 | 1.00 | 0.84 | U | <1.00 | 1.00 | |
| 2-Butanone | 78-93-3 | ~ | 4.7 | D | 1.3 | D | 2.9 | D | 5.01 | 1.00 | 3.45 | 1.00 | 7.66 | 1.00 | 19.4 | 1.00 | 28.2 | 1.00 | 1.09 | 1.00 | 2.3 | D | 10.2 | 12 | D | <1.00 | 1.00 | 0.69 | D | <1.00 | 1.00 | 0.69 | D | <1.00 | 1.00 |
| 2-Hexanone | 591-78-6 | ~ | 0.8 | J | 0.78 | U | 0.85 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.9 | U | <1.00 | 1.00 | 0.77 | U | <1.00 | 1.00 | 0.96 | U | <1.00 | 1.00 | |
| 3-Chloropropene | 107-05-1 | ~ | 1.5 | U | 1.5 | U | 1.6 | U | NT | | NT | | NT | | NT | | NT | | NT | | NT | | 1.7 | U | NT | | 1.5 | U | NT | | 1.8 | U | NT | | |
| 4-Methyl-2-pentanone | 108-10-1 | ~ | 1.6 | D | 0.39 | U | 0.42 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.45 | U | <1.00 | 1.00 | 0.38 | U | <1.00 | 1.00 | 0.48 | U | <1.00 | 1.00 | |
| Acetone | 67-64-1 | ~ | 98 | D | 10 | D | 12 | D | 73.6 | 1.00 | 106 | 1.00 | 44.2 | 1.00 | 79.5 | 1.00 | 140 | 1.00 | 33 | 1.00 | 3300 | D | 13.4 | 100 | D | 10.2 | 100 | D | 7.2 | D | 13.5 | 100 | D | 13.5 | 100 |
| Acrylonitrile | 107-13-1 | ~ | 0.21 | U | 0.21 | U | 0.23 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.24 | U | <1.00 | 1.00 | 0.2 | U | <1.00 | 1.00 | 0.25 | U | <1.00 | 1.00 | |
| Benzene | 71-43-2 | ~ | 1.3 | D | 0.48 | D | 1.1 | D | 4.09 | 1.00 | 5.49 | 1.00 | 2.5 | 1.00 | 3.35 | 1.00 | 5.56 | 1.00 | 1.02 | 1.00 | 0.63 | D | <1.00 | 1.00 | 0.69 | D | <1.00 | 1.00 | 0.56 | D | 1.05 | 1.00 | | | |
| Benzyl chloride | 100-44-7 | ~ | 0.5 | U | 0.49 | U | 0.54 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.57 | U | <1.00 | 1.00 | 0.48 | U | <1.00 | 1.00 | 0.6 | U | <1.00 | 1.00 | |
| Bromodichloromethane | 75-27-4 | ~ | 0.65 | U | 0.63 | U | 0.69 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 0.73 | U | <1.00 | 1.00 | 0.63 | U | <1.00 | 1.00 | 0.78 | U | <1.00 | 1.00 | |
| Bromofom | 75-25-2 | ~ | 1 | U | 0.98 | U | 1.1 | U | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | <1.00 | 1.00 | 1.1 | U | <1.00 | 1.00 | 0.97 | U | <1.00 | 1.00 | 1.2 | U | <1.00 | 1.00 | |
| Bromomethane | 74-83-9 | ~ | 0.38 | U</ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |

Table 3
Volatile Organic Compounds in Groundwater (ug/L)
EPA Method 8260
NYSDEC BCP No. C203119

| Sample ID | Park ID | Sampling Date | Client Matrix | Compound | CAS Number | MW-4 | | MW-6 | | MW-20 | | MW-22 | | MW-23 | | MW-26 | | MW-28 | | SP-1 | | DW-1 | | DW-2 | |
|---|-------------|---------------|---------------|----------|------------|------------|--------|------------|--------|------------|---------|------------|---------|------------|---------|---------|------------|---------|------------|------------|---------|------------|---------|------------|--------|
| | | | | | | 21A0155-01 | 1/5/21 | 21A0155-02 | 1/5/21 | 23F1735-01 | 6/27/23 | 22K0882-01 | 5/16/22 | 22K0882-02 | 5/16/22 | CPS3375 | 11/16/2023 | CPS3376 | 11/16/2023 | 22D1276-01 | 4/25/22 | 23C0637-01 | 4/25/22 | 23C0637-02 | 3/9/23 |
| Volatile Organics, NIDEP/TCL/Part 375 List | | | | | | | | | | | | | | | | | | | | | | | | | |
| Dilution Factor | | | | | | | | | | | | | | | | | | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | 630-20-6 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.216 | U | 0.216 | U |
| 1,1,1-Trichloroethane | 71-55-6 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | <1.0 | 10 | <5.0 | 5 | 40 | U | 0.266 | U | 0.266 | U |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.256 | U | 0.256 | U |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 76-13-1 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.286 | U | 0.286 | U |
| 1,1,2-Trichloroethane | 79-00-5 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.249 | U | 0.249 | U |
| 1,1-Dichloroethane | 75-34-3 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | <1.0 | 10 | <5.0 | 5 | 40 | U | 0.272 | U | 0.272 | U |
| 1,1-Dichloroethylene | 75-35-4 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | NT | 5.99 | U | 70.7 | U | 0.75 | U | <1.0 | 1 | 40 | U | 0.327 | U | 61.8 | U | |
| 1,2,3-Trichlorobenzene | 87-61-6 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.222 | U | 0.222 | U | |
| 1,2,3-Trichloropropane | 96-18-4 | 0.04 | 1 | U | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.273 | U | 0.273 | U | |
| 1,2,4-Trichlorobenzene | 120-82-1 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | NT | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.138 | U | 0.138 | U | |
| 1,2,4-Trimethylbenzene | 95-63-6 | 5 | 5.4 | D | 0.24 | J | NT | 0.2 | U | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.31 | U | 0.31 | U | 0.31 | U |
| 1,2-Dibromo-3-chloropropane | 96-12-8 | 0.04 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.432 | U | 0.432 | U | 0.432 | U |
| 1,2-Dibromoethane | 106-93-4 | 0.0006 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.215 | U | 0.215 | U | 0.215 | U |
| 1,2-Dichlorobenzene | 95-50-1 | 3 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <1.0 | 2 | <1.0 | 1 | 40 | U | 0.27 | U | 0.27 | U | 0.27 | U |
| 1,2-Dichloroethane | 107-06-2 | 0.6 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <1.2 | 1.2 | <0.60 | 0.6 | 40 | U | 0.377 | U | 0.377 | U | 0.377 | U |
| 1,2-Dichloropropane | 78-87-5 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.327 | U | 0.327 | U | 0.327 | U |
| 1,3,5-Trimethylbenzene | 108-67-8 | 5 | 1 | JD | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.347 | U | 0.347 | U | 0.347 | U |
| 1,3-Dichlorobenzene | 541-73-1 | 3 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.283 | U | 0.283 | U | 0.283 | U |
| 1,4-Dichlorobenzene | 106-46-7 | 3 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.36 | J | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.311 | U | 0.311 | U | 0.311 | U | 0.311 | U |
| 1,4-Dioxane | 123-91-1 | ~ | 200 | U | 40 | U | NT | 40 | U | 40 | U | 40 | U | NT | U | NT | U | 8000 | U | 35.3 | U | 35.3 | U | 35.3 | U |
| 2-Butanone | 78-93-3 | 50 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <5.0 | 5 | <2.5 | 2.5 | 40 | U | 11.3 | U | 11.3 | U | 11.3 | U |
| 2-Hexanone | 591-78-6 | 50 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <5.0 | 5 | <2.5 | 2.5 | 40 | U | 0.32 | U | 0.32 | U | 0.32 | U |
| 4-Methyl-2-pentanone | 108-10-1 | ~ | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <5.0 | 5 | <2.5 | 2.5 | 40 | U | 0.365 | U | 0.365 | U | 0.365 | U |
| Acetone | 67-64-1 | 50 | 5 | U | 2.3 | U | NT | 1 | U | 1 | U | 1 | U | <1.0 | 10 | 19 | U | 200 | U | 9.72 | U | 9.72 | U | 9.72 | U |
| Acrolein | 107-02-8 | ~ | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <1.0 | 10 | <5.0 | 5 | 40 | U | 0.447 | U | 0.447 | U | 0.447 | U |
| Acrylonitrile | 107-13-1 | ~ | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <1.0 | 10 | <5.0 | 5 | 40 | U | 0.422 | U | 0.422 | U | 0.422 | U |
| Benzene | 71-43-2 | 1 | 1 | JD | 0.2 | U | NT | 1.04 | U | 0.7 | U | <1.4 | 1.4 | <0.70 | 0.7 | 40 | U | 0.279 | U | 1.32 | U | 1.32 | U | 1.32 | U |
| Bromochloromethane | 74-97-5 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.354 | U | 0.354 | U | 0.354 | U |
| Bromodichloromethane | 75-27-4 | 50 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.245 | U | 0.245 | U | 0.245 | U |
| Bromofrom | 75-25-2 | 50 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <2.0 | 10 | <1.0 | 1 | 40 | U | 0.163 | U | 0.163 | U | 0.163 | U |
| Bromomethane | 74-83-9 | ~ | 1.2 | JD | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <1.0 | 10 | <5.0 | 5 | 40 | U | 0.119 | U | 0.119 | U | 0.119 | U |
| Carbon disulfide | 75-15-0 | ~ | 1 | U | 0.23 | J | NT | 0.23 | J | 0.26 | J | <2.0 | 2 | 1.9 | U | 1 | 40 | U | 0.362 | U | 0.362 | U | 0.362 | U | |
| Carbon tetrachloride | 56-23-5 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.204 | U | 0.204 | U | 0.204 | U |
| Chlorobenzene | 108-90-7 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <1.0 | 10 | <5.0 | 5 | 40 | U | 0.284 | U | 0.284 | U | 0.284 | U |
| Chloroethane | 75-00-3 | 5 | 1 | U | 0.2 | U | NT | 23.3 | U | 0.2 | U | 0.2 | U | <1.0 | 10 | <5.0 | 5 | 40 | U | 0.448 | U | 9.22 | U | 9.22 | U |
| Chloroform | 67-66-3 | 7 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | 1.1 | 10 | 0.35 | U | 40 | U | 2.9 | U | 2.9 | U | 2.9 | U |
| Chloromethane | 74-87-3 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <1.0 | 10 | <5.0 | 5 | 40 | U | 0.372 | U | 0.372 | U | 0.372 | U |
| cis-1,2-Dichloroethylene | 156-59-2 | 5 | 2100 | D | 25 | U | NT | 4620 | D | 31000 | D | 600 | D | 20 | <1.0 | 1 | 40 | U | 10000 | D | 1.37 | U | 22100 | D | |
| cis-1,3-Dichloropropylene | 10061-01-5 | 0.4 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | <0.80 | 0.8 | <0.40 | 0.4 | 40 | U | 0.262 | U | 0.262 | U | 0.262 | U | 0.262 | U |
| Cyclohexane | 110-82-7 | ~ | 5.4 | D | 4.1 | U | NT | 0.53 | U | 4.19 | U | NT | U | NT | U | NT | U | 40 | U | 0.491 | U | 2.24 | U | 2.24 | U |
| Dibromochloromethane | 124-48-1 | 50 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.146 | U | 0.146 | U | 0.146 | U |
| Dibromomethane | 74-95-3 | ~ | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.203 | U | 0.203 | U | 0.203 | U |
| Dichlorodifluoromethane | 75-71-8 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.451 | U | 0.451 | U | 0.451 | U |
| Ethyl Benzene | 100-41-4 | 5 | 2 | JD | 0.21 | J | NT | 0.21 | J | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.29 | U | 1.17 | U | 1.17 | U | 1.17 | U |
| Hexachlorobutadiene | 87-68-3 | 0.5 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | <1.0 | 1 | <0.50 | 0.5 | 40 | U | 0.241 | U | 0.241 | U | 0.241 | U |
| Isopropylbenzene | 98-82-8 | 5 | 16 | D | 4.9 | U | NT | 0.2 | U | 4.74 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.405 | U | 0.405 | U | 0.405 | U | 0.405 | U |
| Methyl acetate | 79-20-9 | ~ | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | U | 0.2 | U | NT | U | NT | U | 40 | U | 0.442 | U | 0.442 | U | 0.442 | U |
| Methyl tert-butyl ether (MTBE) | 1634-04-4 | 10 | 1 | U | 0.2 | U | NT | 0.35 | J | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.33 | J | 0.244 | U | 0.244 | U | 0.244 | U |
| Methylcyclohexane | 108-87-2 | ~ | 24 | D | 8 | U | NT | 0.2 | U | 6.81 | U | NT | U | 0.477 | U | 0.477 | U | 40 | U | 0.477 | U | 2.77 | U | 2.77 | U |
| Methylene chloride | 75-09-2 | 5 | 5 | D | 1 | U | NT | 1.14 | J | 1 | U | <6.0 | 6 | <3.0 | 3 | 40 | U | 0.397 | U | 0.397 | U | 0.397 | U | 0.397 | U |
| n-Butylbenzene | 104-51-8 | 5 | 8.8 | D | 3 | U | NT | 0.2 | U | 0.72 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.399 | U | 0.399 | U | 0.399 | U | 0.399 | U |
| n-Propylbenzene | 103-65-1 | 5 | 3.7 | D | 2.5 | U | NT | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.384 | U | 0.384 | U | 0.384 | U | 0.384 | U |
| p-Xylene | 95-47-6 | 5 | 1.7 | JD | 0.2 | U | NT | 0.2 | U | 5.31 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.261 | U | 0.261 | U | 0.261 | U | 0.261 | U |
| p- & m- Xylenes | 179601-23-1 | 5 | 2.6 | JD | 0.5 | U | NT | 0.5 | U | 1.23 | U | <2.0 | 2 | <1.0 | 1 | 100 | U | 0.578 | U | 0.578 | U | 0.578 | U | 0.578 | U |
| p-Isopropyltoluene | 99-87-6 | 5 | 5.6 | D | 0.33 | J | NT | 0.2 | U | 0.2 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.377 | U | 0.377 | U | 0.377 | U | 0.377 | U |
| sec-Butylbenzene | 135-98-8 | 5 | 16 | D | 8.4 | U | NT | 0.2 | U | 3.75 | U | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.444 | U | 0.444 | U | 0.444 | U | 0.444 | U |
| Styrene | 100-42-5 | 5 | 1 | U | 0.2 | U | NT | 0.2 | U | 0.2 | J | <2.0 | 2 | <1.0 | 1 | 40 | U | 0.255 | U | 0.255 | U | 0.255 | U | 0.255 | U |
| tert-Butyl alcohol (TBA) | 75-65-0 | ~ | 2.5 | U | 0.5 | U | NT | 0.5 | U | 0.5 | U | NT | U | 1.17 | U | 1.17 | U | 100 | U | 0.608 | U | 0.608 | U | 0.608 | U |
| tert-Butylbenzene | 98-06-6 | 5 | 1.1 | JD | 0.77 | U | NT | 0.2 | U | 0.44 | J | <2.0 | 2</ | | | | | | | | | | | | |

Soil Vapor/Indoor Air Matrix A

May 2017

Analytes Assigned:

Trichloroethene (TCE), *cis*-1,2-Dichloroethene (*c*-1,2-DCE), 1,1-Dichloroethene (1,1-DCE), Carbon Tetrachloride

| SUB-SLAB VAPOR CONCENTRATION of COMPOUND (mcg/m ³) | INDOOR AIR CONCENTRATION of COMPOUND (mcg/m ³) | | |
|--|--|----------------------|---|
| | < 0.2 | 0.2 to < 1 | 1 and above |
| < 6 | 1. No further action | 2. No Further Action | 3. IDENTIFY SOURCE(S) and RESAMPLE or MITIGATE |
| 6 to < 60 | 4. No further action | 5. MONITOR | 6. MITIGATE |
| 60 and above | 7. MITIGATE | 8. MITIGATE | 9. MITIGATE |

mcg/m³ = micrograms per cubic meter

No further action: No additional actions are recommended to address human exposures.

Identify Source(s) and Resample or Mitigate: We recommend that reasonable and practical actions be taken to identify the source(s) affecting the indoor air quality and that actions be implemented to reduce indoor air concentrations to within background ranges. For example, if an indoor or outdoor air source is identified, we recommend the appropriate party implement actions to reduce the levels. In the event that indoor or outdoor sources are not readily identified or confirmed, resampling (which might include additional sub-slab vapor and indoor air sampling locations) is recommended to demonstrate that SVI mitigation actions are not needed. Based on the information available, mitigation might also be recommended when soil vapor intrusion cannot be ruled out.

Monitor: We recommend monitoring (sampling on a recurring basis), including but not necessarily limited to sub-slab vapor, basement air and outdoor air sampling, to determine whether concentrations in the indoor air or sub-slab vapor have changed and/or to evaluate temporal influences. Monitoring might also be recommended to determine whether existing building conditions (e.g., positive pressure heating, ventilation and air-conditioning systems) are maintaining the desired mitigation endpoint and to determine whether changes are needed. The type and frequency of monitoring is determined based on site-, building- and analyte-specific information, taking into account applicable environmental data and building operating conditions. Monitoring is an interim measure required to evaluate exposures related to soil vapor intrusion until contaminated environmental media are remediated.

Mitigate: We recommend mitigation to minimize current or potential exposures associated with soil vapor intrusion. The most common mitigation methods are sealing preferential pathways in conjunction with installing a sub-slab depressurization system and changing the pressurization of the building in conjunction with monitoring. The type, or combination of types, of mitigation is determined on a building-specific basis, taking into account building construction and operating conditions. Mitigation is considered a temporary measure implemented to address exposures related to soil vapor intrusion until contaminated environmental media are remediated.

These general recommendations are made with consideration being given to the additional notes on page 2.

ADDITIONAL NOTES FOR MATRIX A

This matrix summarizes actions recommended to address current and potential exposures related to soil vapor intrusion. To use the matrix appropriately as a tool in the decision-making process, the following should be noted:

- [1] The matrix is generic. As such, it may be appropriate to modify a recommended action to accommodate analyte-specific, building-specific conditions (e.g., dirt floor in basement, crawl spaces, thick slabs, current occupancy, etc.), and/or factors provided in Section 3.2 of the guidance (e.g., current land use, environmental conditions, etc.). For example, collection of additional samples may be recommended when the matrix indicates "no further action" for a particular building, but the results of adjacent buildings (especially sub-slab vapor results) indicate a need to take actions to address exposures related to soil vapor intrusion. Mitigation might be recommended when the results of multiple contaminants indicate monitoring is recommended. Proactive actions may be proposed at any time. For example, the party implementing the actions may decide to install sub-slab depressurization systems on buildings where the matrix indicates "no further action" or "monitoring." Such an action might be undertaken for reasons other than public health (e.g., seeking community acceptance, reducing costs, etc.). However, actions implemented *in lieu* of sampling will typically be expected to be captured in the final engineering report and site management plan, and might not rule out the need for post-implementation sampling (e.g., to document effectiveness or to support terminating the action).
- [2] Actions provided in the matrix are specific to addressing human exposures. Implementation of these actions does not preclude investigating possible sources of soil vapor contamination, nor does it preclude remediating contaminated soil vapor or the source of soil vapor contamination.
- [3] Appropriate care should be taken during all aspects of sample collection to ensure that high quality data are obtained. Since the data are being used in the decision-making process, the laboratory analyzing the environmental samples must have current Environmental Laboratory Approval Program (ELAP) certification for the appropriate analyte and environmental matrix combinations. Furthermore, samples should be analyzed by methods that can achieve a minimum reporting limit of 0.20 microgram per cubic meter for indoor and outdoor air samples. For sub-slab vapor samples and dirt floor soil vapor samples, a minimum reporting limit of 1 microgram per cubic meter is recommended.
- [4] Sub-slab vapor and indoor air samples are typically collected when the likelihood of soil vapor intrusion is considered to be the greatest (i.e., worst-case conditions). If samples are collected at other times (typically, samples collected outside of the heating season), then resampling during worst-case conditions might be appropriate to verify that actions taken to address exposures related to soil vapor intrusion are protective of human health.
- [5] When current exposures are attributed to sources other than soil vapor intrusion, the agencies should be given documentation (e.g., applicable environmental data, completed indoor air sampling questionnaire, digital photographs, etc.) to support a proposed action other than that provided in the matrix box and to support agency assessment and follow-up.
- [6] The party responsible for implementing the recommended actions will differ depending upon several factors, including but not limited to the following: the identified source of the volatile chemicals, the environmental remediation program, and analyte-specific, site-specific and building-specific factors.

Soil Vapor/Indoor Air Matrix B

May 2017

Analytes Assigned:

Tetrachloroethene (PCE), 1,1,1-Trichloroethane (1,1,1-TCA), Methylene Chloride

| SUB-SLAB VAPOR CONCENTRATION of COMPOUND (mcg/m ³) | INDOOR AIR CONCENTRATION of COMPOUND (mcg/m ³) | | |
|--|--|----------------------|---|
| | < 3 | 3 to < 10 | 10 and above |
| < 100 | 1. No further action | 2. No Further Action | 3. IDENTIFY SOURCE(S) and RESAMPLE or MITIGATE |
| 100 to < 1,000 | 4. No further action | 5. MONITOR | 6. MITIGATE |
| 1,000 and above | 7. MITIGATE | 8. MITIGATE | 9. MITIGATE |

mcg/m³ = micrograms per cubic meter

No further action: No additional actions are recommended to address human exposures.

Identify Source(s) and Resample or Mitigate: We recommend that reasonable and practical actions be taken to identify the source(s) affecting the indoor air quality and that actions be implemented to reduce indoor air concentrations to within background ranges. For example, if an indoor or outdoor air source is identified, we recommend the appropriate party implement actions to reduce the levels. In the event that indoor or outdoor sources are not readily identified or confirmed, resampling (which might include additional sub-slab vapor and indoor air sampling locations) is recommended to demonstrate that SVI mitigation actions are not needed. Based on the information available, mitigation might also be recommended when soil vapor intrusion cannot be ruled out.

Monitor: We recommend monitoring (sampling on a recurring basis), including but not necessarily limited to sub-slab vapor, basement air and outdoor air sampling, to determine whether concentrations in the indoor air or sub-slab vapor have changed and/or to evaluate temporal influences. Monitoring might also be recommended to determine whether existing building conditions (e.g., positive pressure heating, ventilation and air-conditioning systems) are maintaining the desired mitigation endpoint and to determine whether changes are needed. The type and frequency of monitoring is determined based on site-, building- and analyte-specific information, taking into account applicable environmental data and building operating conditions. Monitoring is an interim measure required to evaluate exposures related to soil vapor intrusion until contaminated environmental media are remediated.

Mitigate: We recommend mitigation to minimize current or potential exposures associated with soil vapor intrusion. The most common mitigation methods are sealing preferential pathways in conjunction with installing a sub-slab depressurization system and changing the pressurization of the building in conjunction with monitoring. The type, or combination of types, of mitigation is determined on a building-specific basis, taking into account building construction and operating conditions. Mitigation is considered a temporary measure implemented to address exposures related to soil vapor intrusion until contaminated environmental media are remediated.

These general recommendations are made with consideration being given to the additional notes on page 2.

ADDITIONAL NOTES FOR MATRIX B

This matrix summarizes actions recommended to address current and potential exposures related to soil vapor intrusion. To use the matrix appropriately as a tool in the decision-making process, the following should be noted:

- [1] The matrix is generic. As such, it may be appropriate to modify a recommended action to accommodate analyte-specific, building-specific conditions (e.g., dirt floor in basement, crawl spaces, thick slabs, current occupancy, etc.), and/or factors provided in Section 3.2 of the guidance (e.g., current land use, environmental conditions, etc.). For example, collection of additional samples may be recommended when the matrix indicates "no further action" for a particular building, but the results of adjacent buildings (especially sub-slab vapor results) indicate a need to take actions to address exposures related to soil vapor intrusion. Mitigation might be recommended when the results of multiple contaminants indicate monitoring is recommended. Proactive actions may be proposed at any time. For example, the party implementing the actions may decide to install sub-slab depressurization systems on buildings where the matrix indicates "no further action" or "monitoring." Such an action might be undertaken for reasons other than public health (e.g., seeking community acceptance, reducing costs, etc.). However, actions implemented *in lieu* of sampling will typically be expected to be captured in the final engineering report and site management plan, and might not rule out the need for post-implementation sampling (e.g., to document effectiveness or to support terminating the action).
- [2] Actions provided in the matrix are specific to addressing human exposures. Implementation of these actions does not preclude investigating possible sources of soil vapor contamination, nor does it preclude remediating contaminated soil vapor or the source of soil vapor contamination.
- [3] Appropriate care should be taken during all aspects of sample collection to ensure that high quality data are obtained. Since the data are being used in the decision-making process, the laboratory analyzing the environmental samples must have current Environmental Laboratory Approval Program (ELAP) certification for the appropriate analyte and environmental matrix combinations. Furthermore, samples should be analyzed by methods that can achieve a minimum reporting limit of 1 microgram per cubic meter for indoor and outdoor air samples. For sub-slab vapor samples and dirt floor soil vapor samples, a minimum reporting limit of 1 microgram per cubic meter is recommended.
- [4] Sub-slab vapor and indoor air samples are typically collected when the likelihood of soil vapor intrusion to occur is considered to be the greatest (i.e., worst-case conditions). If samples are collected at other times (typically, samples collected outside of the heating season), then resampling during worst-case conditions might be appropriate to verify that actions taken to address exposures related to soil vapor intrusion are protective of human health.
- [5] When current exposures are attributed to sources other than soil vapor intrusion, the agencies should be given documentation (e.g., applicable environmental data, completed indoor air sampling questionnaire, digital photographs, etc.) to support a proposed action other than that provided in the matrix box and to support agency assessment and follow-up.
- [6] The party responsible for implementing the recommended actions will differ depending upon several factors, including but not limited to the following: the identified source of the volatile chemicals, the environmental remediation program, and analyte-specific, site-specific and building-specific factors.

Soil Vapor/Indoor Air Matrix C

May 2017

Analytes Assigned:

Vinyl Chloride

| | | INDOOR AIR CONCENTRATION of COMPOUND (mcg/m ³) | |
|--|----------------------|--|--|
| SUB-SLAB VAPOR CONCENTRATION of COMPOUND (mcg/m ³) | | | |
| | < 0.2 | 0.2 and above | |
| < 6 | 1. No further action | 3. IDENTIFY SOURCE(S) and RESAMPLE or MITIGATE | |
| 6 to < 60 | 4. MONITOR | 6. MITIGATE | |
| 60 and above | 7. MITIGATE | 9. MITIGATE | |

mcg/m³ = micrograms per cubic meter

No further action: No additional actions are recommended to address human exposures.

Identify Source(s) and Resample or Mitigate: We recommend that reasonable and practical actions be taken to identify the source(s) affecting the indoor air quality and that actions be implemented to reduce indoor air concentrations to within background ranges. For example, if an indoor or outdoor air source is identified, we recommend the appropriate party implement actions to reduce the levels. In the event that indoor or outdoor sources are not readily identified or confirmed, resampling (which might include additional sub-slab vapor and indoor air sampling locations) is recommended to demonstrate that SVI mitigation actions are not needed. Based on the information available, mitigation might also be recommended when soil vapor intrusion cannot be ruled out.

Monitor: We recommend monitoring (sampling on a recurring basis), including but not necessarily limited to sub-slab vapor, basement air and outdoor air sampling, to determine whether concentrations in the indoor air or sub-slab vapor have changed and/or to evaluate temporal influences. Monitoring might also be recommended to determine whether existing building conditions (e.g., positive pressure heating, ventilation and air-conditioning systems) are maintaining the desired mitigation endpoint and to determine whether changes are needed. The type and frequency of monitoring is determined based on site-, building- and analyte-specific information, taking into account applicable environmental data and building operating conditions. Monitoring is an interim measure required to evaluate exposures related to soil vapor intrusion until contaminated environmental media are remediated.

Mitigate: We recommend mitigation to minimize current or potential exposures associated with soil vapor intrusion. The most common mitigation methods are sealing preferential pathways in conjunction with installing a sub-slab depressurization system and changing the pressurization of the building in conjunction with monitoring. The type, or combination of types, of mitigation is determined on a building-specific basis, taking into account building construction and operating conditions. Mitigation is considered a temporary measure implemented to address exposures related to soil vapor intrusion until contaminated environmental media are remediated.

These general recommendations are made with consideration being given to the additional notes on page 2.

ADDITIONAL NOTES FOR MATRIX C

This matrix summarizes actions recommended to address current and potential exposures related to soil vapor intrusion. To use the matrix appropriately as a tool in the decision-making process, the following should be noted:

- [1] The matrix is generic. As such, it may be appropriate to modify a recommended action to accommodate analyte-specific, building-specific conditions (e.g., dirt floor in basement, crawl spaces, thick slabs, current occupancy, etc.), and/or factors provided in Section 3.2 of the guidance (e.g., current land use, environmental conditions, etc.). For example, collection of additional samples may be recommended when the matrix indicates "no further action" for a particular building, but the results of adjacent buildings (especially sub-slab vapor results) indicate a need to take actions to address exposures related to soil vapor intrusion. Mitigation might be recommended when the results of multiple contaminants indicate monitoring is recommended. Proactive actions may be proposed at any time. For example, the party implementing the actions may decide to install sub-slab depressurization systems on buildings where the matrix indicates "no further action" or "monitoring." Such an action might be undertaken for reasons other than public health (e.g., seeking community acceptance, reducing costs, etc.). However, actions implemented *in lieu* of sampling will typically be expected to be captured in the final engineering report and site management plan, and might not rule out the need for post-implementation sampling (e.g., to document effectiveness or to support terminating the action).
- [2] Actions provided in the matrix are specific to addressing human exposures. Implementation of these actions does not preclude investigating possible sources of soil vapor contamination, nor does it preclude remediating contaminated soil vapor or the source of soil vapor contamination.
- [3] Appropriate care should be taken during all aspects of sample collection to ensure that high quality data are obtained. Since the data are being used in the decision-making process, the laboratory analyzing the environmental samples must have current Environmental Laboratory Approval Program (ELAP) certification for the appropriate analyte and environmental matrix combinations. Furthermore, samples should be analyzed by methods that can achieve a minimum reporting limit of 0.20 microgram per cubic meter for indoor and outdoor air samples. For sub-slab vapor samples and dirt floor soil vapor samples, a minimum reporting limit of 1 microgram per cubic meter is recommended.
- [4] Sub-slab vapor and indoor air samples are typically collected when the likelihood of soil vapor intrusion is considered to be the greatest (i.e., worst-case conditions). If samples are collected at other times (typically, samples collected outside of the heating season), then resampling during worst-case conditions might be appropriate to verify that actions taken to address exposures related to soil vapor intrusion are protective of human health.
- [5] When current exposures are attributed to sources other than soil vapor intrusion, the agencies should be given documentation (e.g., applicable environmental data, completed indoor air sampling questionnaire, digital photographs, etc.) to support a proposed action other than that provided in the matrix box and to support agency assessment and follow-up.
- [6] The party responsible for implementing the recommended actions will differ depending upon several factors, including but not limited to the following: the identified source of the volatile chemicals, the environmental remediation program, and analyte-specific, site-specific and building-specific factors.

Soil Vapor/Indoor Air Matrix D

February 2024

Analytes Assigned:

Benzene, ethylbenzene, naphthalene, cyclohexane, isooctane (2,2,4-trimethylpentane), 1,2,4-trimethylbenzene, 1,3,5-trimethylbenzene, o-xylene

| SUB-SLAB VAPOR CONCENTRATION of COMPOUND (mcg/m ³) | INDOOR AIR CONCENTRATION of COMPOUND (mcg/m ³) | | |
|--|--|----------------------|---|
| | < 2 | 2 to < 10 | 10 and above |
| < 60 | 1. No further action | 2. No Further Action | 3. IDENTIFY SOURCE(S) or RESAMPLE or MITIGATE |
| 60 to < 600 | 4. No further action | 5. MONITOR | 6. MITIGATE |
| 600 and above | 7. MITIGATE | 8. MITIGATE | 9. MITIGATE |

mcg/m³ = micrograms per cubic meter

No further action: No additional actions are recommended to address human exposures.

Identify Source(s) or Resample or Mitigate: We recommend that reasonable and practical actions be taken to identify the source(s) affecting the indoor air quality and that actions be implemented to reduce indoor air concentrations to within background ranges. For example, if an indoor or outdoor air source is identified, we recommend the appropriate party implement actions to reduce the levels. In the event that indoor or outdoor sources are not readily identified or confirmed, resampling (which might include additional sub-slab vapor and indoor air sampling locations) is recommended to demonstrate that SVI mitigation actions are not needed. Based on the information available, mitigation might also be recommended when soil vapor intrusion cannot be ruled out.

Monitor: We recommend monitoring (sampling on a recurring basis), including but not necessarily limited to sub-slab vapor, basement air and outdoor air sampling, to determine whether concentrations in the indoor air or sub-slab vapor have changed and/or to evaluate temporal influences. Monitoring might also be recommended to determine whether existing building conditions (e.g., positive pressure heating, ventilation, and air-conditioning systems) are maintaining the desired mitigation endpoint and to determine whether changes are needed. The type and frequency of monitoring is determined based on site-, building- and analyte-specific information, taking into account applicable environmental data and building operating conditions. Monitoring is an interim measure required to evaluate exposures related to soil vapor intrusion until contaminated environmental media are remediated.

Mitigate: We recommend mitigation to minimize current or potential exposures associated with soil vapor intrusion. The most common mitigation methods are sealing preferential pathways in conjunction with installing a sub-slab depressurization system and changing the pressurization of the building in conjunction with monitoring. The type, or combination of types, of mitigation is determined on a building -specific basis, taking into account building construction and operating conditions. Mitigation is considered a temporary measure implemented to address exposures related to soil vapor intrusion until contaminated environmental media are remediated.

These general recommendations are made with consideration being given to the additional notes on page 2.

ADDITIONAL NOTES FOR MATRIX D

This matrix summarizes actions recommended to address current and potential exposures related to soil vapor intrusion. To use the matrix appropriately as a tool in the decision-making process, the following should be noted:

- [1] The matrix is generic. As such, it may be appropriate to modify a recommended action to accommodate analyte-specific, building-specific conditions (e.g., dirt floor in basement, crawl spaces, thick slabs, current occupancy, etc.), and/or factors provided in Section 3.2 of the guidance (e.g., current land use, environmental conditions, etc.). For example, collection of additional samples may be recommended when the matrix indicates "no further action" for a particular building, but the results of adjacent buildings (especially sub-slab vapor results) indicate a need to take actions to address exposures related to soil vapor intrusion. Mitigation might be recommended when the results of multiple contaminants indicate monitoring is recommended. Proactive actions may be proposed at any time. For example, the party implementing the actions may decide to install sub-slab depressurization systems on buildings where the matrix indicates "no further action" or "monitoring." Such an action might be undertaken for reasons other than public health (e.g., seeking community acceptance, reducing costs, etc.). However, actions implemented *in lieu* of sampling will typically be expected to be captured in the final engineering report and site management plan, and might not rule out the need for post-implementation sampling (e.g., to document effectiveness or to support terminating the action).
- [2] Actions provided in the matrix are specific to addressing human exposures. Implementation of these actions does not preclude investigating possible sources of soil vapor contamination, nor does it preclude remediating contaminated soil vapor or the source of soil vapor contamination.
- [3] Appropriate care should be taken during all aspects of sample collection to ensure that high quality data are obtained. Since the data are being used in the decision-making process, the laboratory analyzing the environmental samples must have current Environmental Laboratory Approval Program (ELAP) certification for the appropriate analyte and environmental matrix combinations. Furthermore, samples should be analyzed by methods that can achieve a minimum reporting limit of 1 microgram per cubic meter for indoor and outdoor air samples. For sub-slab vapor samples and dirt floor soil vapor samples, a minimum reporting limit of 1 microgram per cubic meter is recommended.
- [4] Sub-slab vapor and indoor air samples are typically collected when the likelihood of soil vapor intrusion to occur is considered to be the greatest (i.e., worst-case conditions). If samples are collected at other times (typically, samples collected outside of the heating season), then resampling during worst-case conditions might be appropriate to verify that actions taken to address exposures related to soil vapor intrusion are protective of human health.
- [5] When current exposures are attributed to sources other than soil vapor intrusion, the agencies should be given documentation (e.g., applicable environmental data, completed indoor air sampling questionnaire, digital photographs, etc.) to support a proposed action other than that provided in the matrix box and to support agency assessment and follow-up.
- [6] The party responsible for implementing the recommended actions will differ depending upon several factors, including but not limited to the following: the identified source of the volatile chemicals, the environmental remediation program, and analyte-specific, site-specific and building-specific factors.

Soil Vapor/Indoor Air Matrix E

February 2024

Analytes Assigned:

m,p-xylene, heptane, hexane

| SUB-SLAB VAPOR CONCENTRATION of COMPOUND (mcg/m ³) | INDOOR AIR CONCENTRATION of COMPOUND (mcg/m ³) | | |
|--|--|----------------------|---|
| | < 6 | 6 to < 20 | 20 and above |
| < 200 | 1. No further action | 2. No Further Action | 3. IDENTIFY SOURCE(S) or RESAMPLE or MITIGATE |
| 200 to < 2,000 | 4. No further action | 5. MONITOR | 6. MITIGATE |
| 2,000 and above | 7. MITIGATE | 8. MITIGATE | 9. MITIGATE |

mcg/m³ = micrograms per cubic meter

No further action: No additional actions are recommended to address human exposures.

Identify Source(s) or Resample or Mitigate: We recommend that reasonable and practical actions be taken to identify the source(s) affecting the indoor air quality and that actions be implemented to reduce indoor air concentrations to within background ranges. For example, if an indoor or outdoor air source is identified, we recommend the appropriate party implement actions to reduce the levels. In the event that indoor or outdoor sources are not readily identified or confirmed, resampling (which might include additional sub-slab vapor and indoor air sampling locations) is recommended to demonstrate that SVI mitigation actions are not needed. Based on the information available, mitigation might also be recommended when soil vapor intrusion cannot be ruled out.

Monitor: We recommend monitoring (sampling on a recurring basis), including but not necessarily limited to sub-slab vapor, basement air and outdoor air sampling, to determine whether concentrations in the indoor air or sub-slab vapor have changed and/or to evaluate temporal influences. Monitoring might also be recommended to determine whether existing building conditions (e.g., positive pressure heating, ventilation, and air-conditioning systems) are maintaining the desired mitigation endpoint and to determine whether changes are needed. The type and frequency of monitoring is determined based on site-, building- and analyte-specific information, taking into account applicable environmental data and building operating conditions. Monitoring is an interim measure required to evaluate exposures related to soil vapor intrusion until contaminated environmental media are remediated.

Mitigate: We recommend mitigation to minimize current or potential exposures associated with soil vapor intrusion. The most common mitigation methods are sealing preferential pathways in conjunction with installing a sub-slab depressurization system and changing the pressurization of the building in conjunction with monitoring. The type, or combination of types, of mitigation is determined on a building -specific basis, taking into account building construction and operating conditions. Mitigation is considered a temporary measure implemented to address exposures related to soil vapor intrusion until contaminated environmental media are remediated.

These general recommendations are made with consideration being given to the additional notes on page 2.

ADDITIONAL NOTES FOR MATRIX E

This matrix summarizes actions recommended to address current and potential exposures related to soil vapor intrusion. To use the matrix appropriately as a tool in the decision-making process, the following should be noted:

- [1] The matrix is generic. As such, it may be appropriate to modify a recommended action to accommodate analyte-specific, building-specific conditions (e.g., dirt floor in basement, crawl spaces, thick slabs, current occupancy, etc.), and/or factors provided in Section 3.2 of the guidance (e.g., current land use, environmental conditions, etc.). For example, collection of additional samples may be recommended when the matrix indicates "no further action" for a particular building, but the results of adjacent buildings (especially sub-slab vapor results) indicate a need to take actions to address exposures related to soil vapor intrusion. Mitigation might be recommended when the results of multiple contaminants indicate monitoring is recommended. Proactive actions may be proposed at any time. For example, the party implementing the actions may decide to install sub-slab depressurization systems on buildings where the matrix indicates "no further action" or "monitoring." Such an action might be undertaken for reasons other than public health (e.g., seeking community acceptance, reducing costs, etc.). However, actions implemented *in lieu* of sampling will typically be expected to be captured in the final engineering report and site management plan, and might not rule out the need for post-implementation sampling (e.g., to document effectiveness or to support terminating the action).
- [2] Actions provided in the matrix are specific to addressing human exposures. Implementation of these actions does not preclude investigating possible sources of soil vapor contamination, nor does it preclude remediating contaminated soil vapor or the source of soil vapor contamination.
- [3] Appropriate care should be taken during all aspects of sample collection to ensure that high quality data are obtained. Since the data are being used in the decision-making process, the laboratory analyzing the environmental samples must have current Environmental Laboratory Approval Program (ELAP) certification for the appropriate analyte and environmental matrix combinations. Furthermore, samples should be analyzed by methods that can achieve a minimum reporting limit of 1 microgram per cubic meter for indoor and outdoor air samples. For sub-slab vapor samples and dirt floor soil vapor samples, a minimum reporting limit of 1 microgram per cubic meter is recommended.
- [4] Sub-slab vapor and indoor air samples are typically collected when the likelihood of soil vapor intrusion to occur is considered to be the greatest (i.e., worst-case conditions). If samples are collected at other times (typically, samples collected outside of the heating season), then resampling during worst-case conditions might be appropriate to verify that actions taken to address exposures related to soil vapor intrusion are protective of human health.
- [5] When current exposures are attributed to sources other than soil vapor intrusion, the agencies should be given documentation (e.g., applicable environmental data, completed indoor air sampling questionnaire, digital photographs, etc.) to support a proposed action other than that provided in the matrix box and to support agency assessment and follow-up.
- [6] The party responsible for implementing the recommended actions will differ depending upon several factors, including but not limited to the following: the identified source of the volatile chemicals, the environmental remediation program, and analyte-specific, site-specific and building-specific factors.

Soil Vapor/Indoor Air Matrix F

February 2024

Analytes Assigned:

Toluene

| SUB-SLAB VAPOR CONCENTRATION of COMPOUND (mcg/m ³) | INDOOR AIR CONCENTRATION of COMPOUND (mcg/m ³) | | |
|--|--|----------------------|---|
| | < 10 | 10 to < 50 | 50 and above |
| < 300 | 1. No Further Action | 2. No Further Action | 3. IDENTIFY SOURCE(S) or RESAMPLE or MITIGATE |
| 300 to < 3,000 | 4. No Further Action | 5. MONITOR | 6. MITIGATE |
| 3,000 and above | 7. MITIGATE | 8. MITIGATE | 9. MITIGATE |

mcg/m³ = micrograms per cubic meter

No further action: No additional actions are recommended to address human exposures.

Identify Source(s) or Resample or Mitigate: We recommend that reasonable and practical actions be taken to identify the source(s) affecting the indoor air quality and that actions be implemented to reduce indoor air concentrations to within background ranges. For example, if an indoor or outdoor air source is identified, we recommend the appropriate party implement actions to reduce the levels. In the event that indoor or outdoor sources are not readily identified or confirmed, resampling (which might include additional sub-slab vapor and indoor air sampling locations) is recommended to demonstrate that SVI mitigation actions are not needed. Based on the information available, mitigation might also be recommended when soil vapor intrusion cannot be ruled out.

Monitor: We recommend monitoring (sampling on a recurring basis), including but not necessarily limited to sub-slab vapor, basement air and outdoor air sampling, to determine whether concentrations in the indoor air or sub-slab vapor have changed and/or to evaluate temporal influences. Monitoring might also be recommended to determine whether existing building conditions (e.g., positive pressure heating, ventilation, and air-conditioning systems) are maintaining the desired mitigation endpoint and to determine whether changes are needed. The type and frequency of monitoring is determined based on site-, building- and analyte-specific information, taking into account applicable environmental data and building operating conditions. Monitoring is an interim measure required to evaluate exposures related to soil vapor intrusion until contaminated environmental media are remediated.

Mitigate: We recommend mitigation to minimize current or potential exposures associated with soil vapor intrusion. The most common mitigation methods are sealing preferential pathways in conjunction with installing a sub-slab depressurization system and changing the pressurization of the building in conjunction with monitoring. The type, or combination of types, of mitigation is determined on a building -specific basis, taking into account building construction and operating conditions. Mitigation is considered a temporary measure implemented to address exposures related to soil vapor intrusion until contaminated environmental media are remediated.

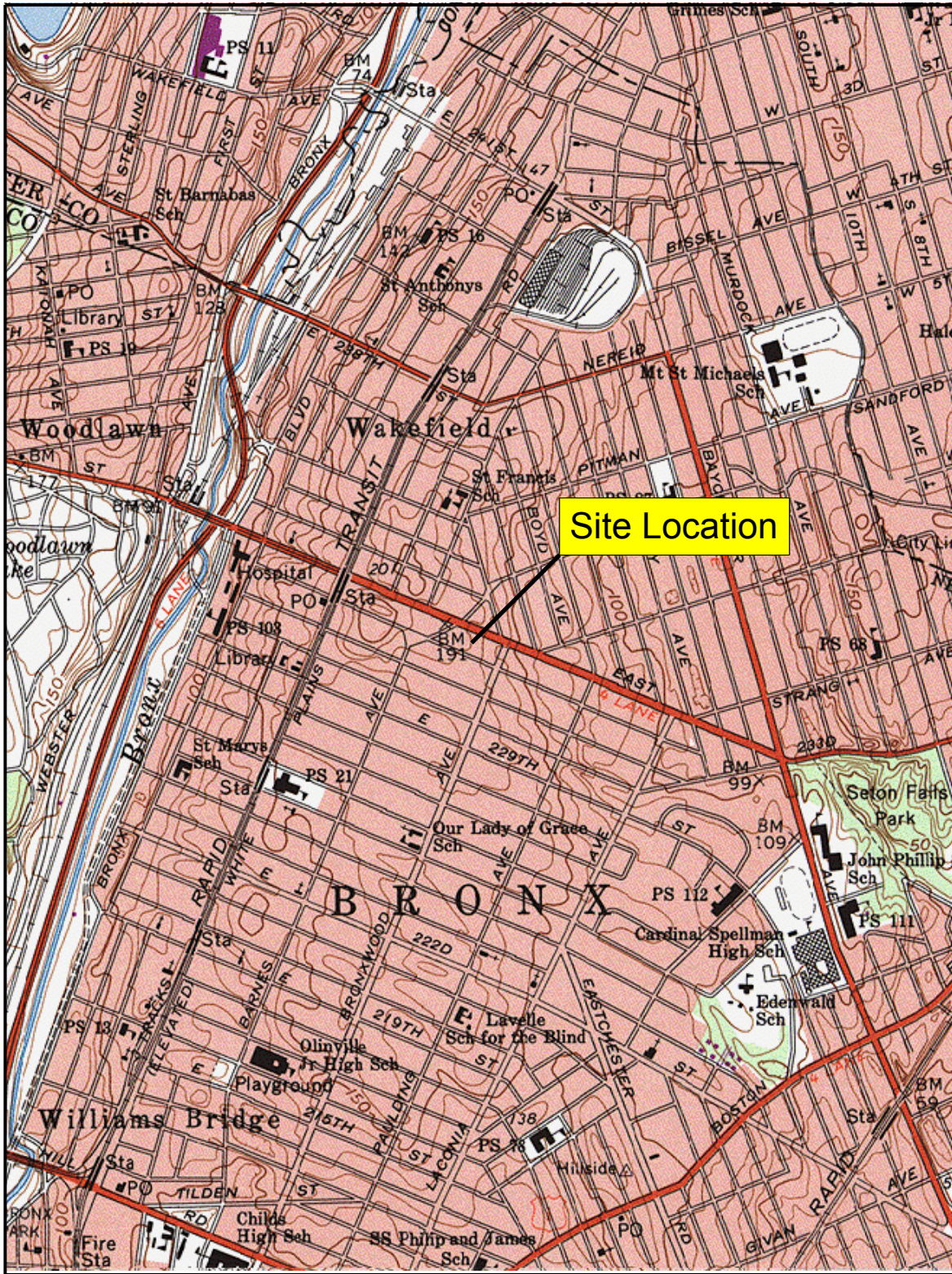
These general recommendations are made with consideration being given to the additional notes on page 2.

ADDITIONAL NOTES FOR MATRIX F

This matrix summarizes actions recommended to address current and potential exposures related to soil vapor intrusion. To use the matrix appropriately as a tool in the decision-making process, the following should be noted:

- [1] The matrix is generic. As such, it may be appropriate to modify a recommended action to accommodate analyte-specific, building-specific conditions (e.g., dirt floor in basement, crawl spaces, thick slabs, current occupancy, etc.), and/or factors provided in Section 3.2 of the guidance (e.g., current land use, environmental conditions, etc.). For example, collection of additional samples may be recommended when the matrix indicates "no further action" for a particular building, but the results of adjacent buildings (especially sub-slab vapor results) indicate a need to take actions to address exposures related to soil vapor intrusion. Mitigation might be recommended when the results of multiple contaminants indicate monitoring is recommended. Proactive actions may be proposed at any time. For example, the party implementing the actions may decide to install sub-slab depressurization systems on buildings where the matrix indicates "no further action" or "monitoring." Such an action might be undertaken for reasons other than public health (e.g., seeking community acceptance, reducing costs, etc.). However, actions implemented *in lieu* of sampling will typically be expected to be captured in the final engineering report and site management plan, and might not rule out the need for post-implementation sampling (e.g., to document effectiveness or to support terminating the action).
- [2] Actions provided in the matrix are specific to addressing human exposures. Implementation of these actions does not preclude investigating possible sources of soil vapor contamination, nor does it preclude remediating contaminated soil vapor or the source of soil vapor contamination.
- [3] Appropriate care should be taken during all aspects of sample collection to ensure that high quality data are obtained. Since the data are being used in the decision-making process, the laboratory analyzing the environmental samples must have current Environmental Laboratory Approval Program (ELAP) certification for the appropriate analyte and environmental matrix combinations. Furthermore, samples should be analyzed by methods that can achieve a minimum reporting limit of 1 microgram per cubic meter for indoor and outdoor air samples. For sub-slab vapor samples and dirt floor soil vapor samples, a minimum reporting limit of 1 microgram per cubic meter is recommended.
- [4] Sub-slab vapor and indoor air samples are typically collected when the likelihood of soil vapor intrusion to occur is considered to be the greatest (i.e., worst-case conditions). If samples are collected at other times (typically, samples collected outside of the heating season), then resampling during worst-case conditions might be appropriate to verify that actions taken to address exposures related to soil vapor intrusion are protective of human health.
- [5] When current exposures are attributed to sources other than soil vapor intrusion, the agencies should be given documentation (e.g., applicable environmental data, completed indoor air sampling questionnaire, digital photographs, etc.) to support a proposed action other than that provided in the matrix box and to support agency assessment and follow-up.
- [6] The party responsible for implementing the recommended actions will differ depending upon several factors, including but not limited to the following: the identified source of the volatile chemicals, the environmental remediation program, and analyte-specific, site-specific and building-specific factors.

Figures



From USGS 7.5 Minute Topographic Map of Mount Vernon, NY Quadrangle



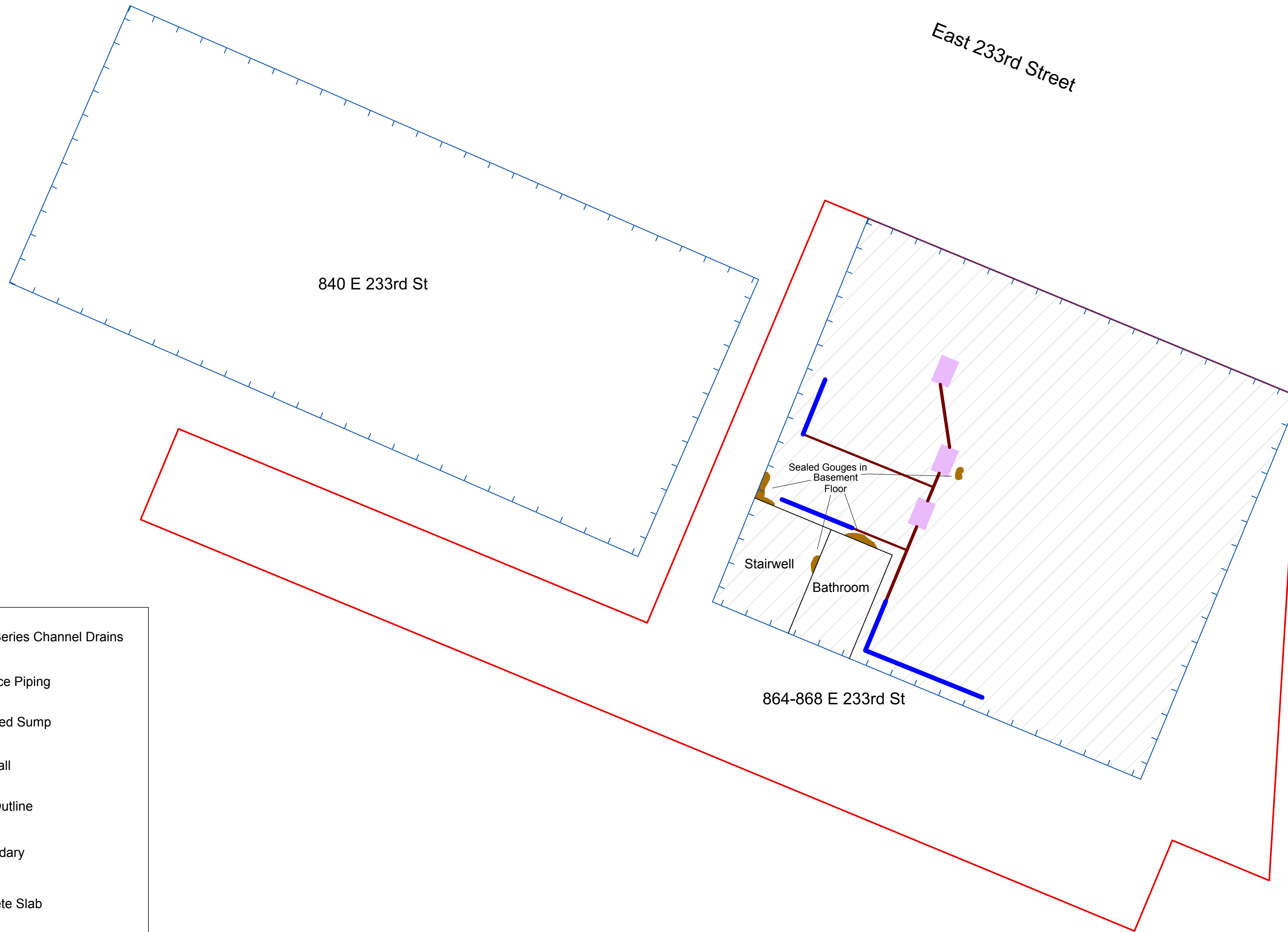
Locational Diagram

Advanced Cleanup Technologies, Inc.
ENVIRONMENTAL CONSULTANTS

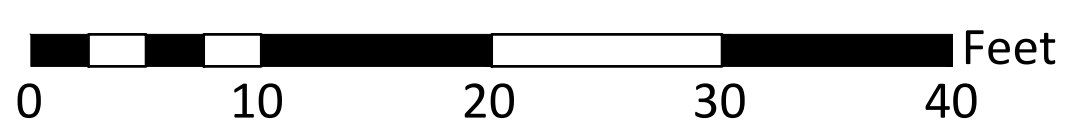
110 Main Street, Suite 103, Port Washington, New York 11050
Tel: 516-441-5800 Fax: 516-441-5511

Project No.: 9628-BXNY Figure No.: 1

Date: 10/18/2018 Scale: Not To Scale



| | |
|--|-------------------------------|
| | NDS Pro Series Channel Drains |
| | Sub-Surface Piping |
| | NDS Sealed Sump |
| | Interior Wall |
| | Building Outline |
| | Site Boundary |
| | 6" Concrete Slab |



Notes

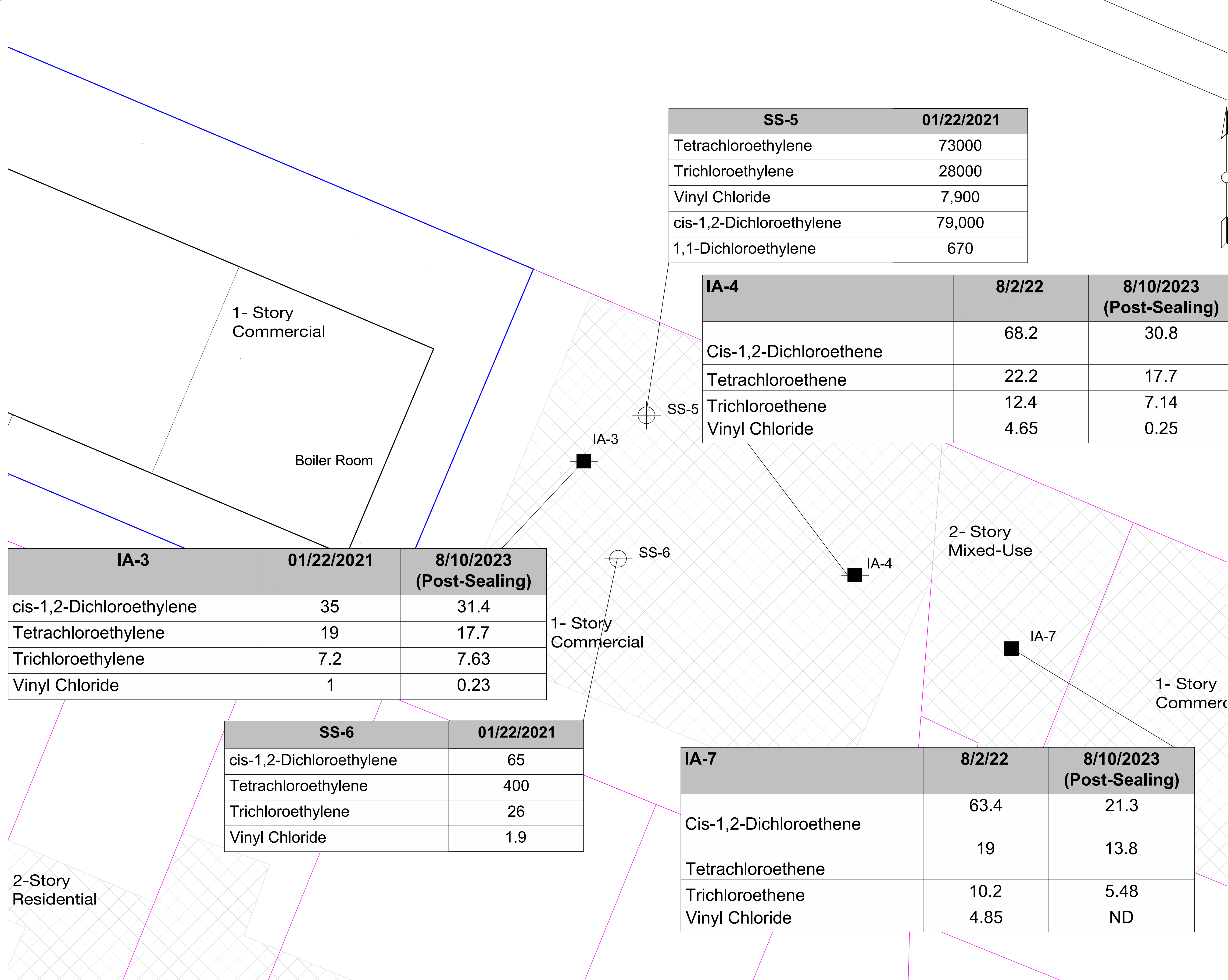
Title
As-Built Site
Diagram

| No. | Revision/Issue | Date |
|-----|----------------|------|
| | | |

Advanced Cleanup Technologies
Environmental Consultants
228 Park Ave S PMB 34864
New York, New York 10003

Project Name and Address
Jay Dry Cleaners
840 E 233rd St
Bronx, NY
10466

| | |
|-----------------------------|--------------------|
| Project BCP Site C203119 | Figure 2 |
| Date 8/21/2024 | |
| Scale As Noted | |



| SS-5 | 01/22/2021 |
|--------------------------|------------|
| Tetrachloroethylene | 73000 |
| Trichloroethylene | 28000 |
| Vinyl Chloride | 7,900 |
| cis-1,2-Dichloroethylene | 79,000 |
| 1,1-Dichloroethylene | 670 |

| IA-4 | 8/2/22 | 8/10/2023 (Post-Sealing) |
|------------------------|--------|-----------------------------|
| Cis-1,2-Dichloroethene | 68.2 | 30.8 |
| Tetrachloroethene | 22.2 | 17.7 |
| Trichloroethene | 12.4 | 7.14 |
| Vinyl Chloride | 4.65 | 0.25 |

| IA-3 | 01/22/2021 | 8/10/2023 (Post-Sealing) |
|--------------------------|------------|-----------------------------|
| cis-1,2-Dichloroethylene | 35 | 31.4 |
| Tetrachloroethylene | 19 | 17.7 |
| Trichloroethylene | 7.2 | 7.63 |
| Vinyl Chloride | 1 | 0.23 |

| SS-6 | 01/22/2021 |
|--------------------------|------------|
| cis-1,2-Dichloroethylene | 65 |
| Tetrachloroethylene | 400 |
| Trichloroethylene | 26 |
| Vinyl Chloride | 1.9 |

| IA-7 | 8/2/22 | 8/10/2023 (Post-Sealing) |
|------------------------|--------|-----------------------------|
| Cis-1,2-Dichloroethene | 63.4 | 21.3 |
| Tetrachloroethene | 19 | 13.8 |
| Trichloroethene | 10.2 | 5.48 |
| Vinyl Chloride | 4.85 | ND |

Notes
Compound unit: ug/m³

| Legend | |
|--------|-----------------------------|
| | BCP Site Boundary |
| | Offsite Site Boundaries |
| | Soil Boring/Monitoring Well |
| | Discrete Water Sample |
| | Soil Boring |
| | Soil Vapor |
| | Outdoor Air |
| | Basement Indoor Air |
| | Floor 1 Indoor Air |
| | IRM |
| | IRM Addendum |

Title
Soil Vapor Intrusion
Spider Diagram

| No. | Revision/Issue | Date |
|-----|----------------|------|
| | | |

Advanced Cleanup Technologies
Environmental Consultants
200 Broadhollow Road-Suite 207
Melville, New York 11747

Project Name and Address
840 East 233rd Street
Bronx, NY 10466

| Project | Figure |
|-----------|----------|
| 9628-BXNY | 3 |
| Date | |
| 3/10/2025 | |
| Scale | As Noted |

Notes

Yellow highlight value exceed the NYSDEC TOGS Standards and Guidance-GA

Compound unit: ug/L
PFAS Unit: ng/L

Legend

| | |
|--|-----------------------------|
| | BCP Site Boundary |
| | Offsite Site Boundaries |
| | Soil Boring/Monitoring Well |
| | Discrete Water Sample |
| | Soil Boring |
| | Soil Vapor |
| | Outdoor Air |
| | Basement Indoor Air |
| | Floor 1 Indoor Air |
| | IRM |
| | IRM Addendum |

Title

Groundwater Exceedance Diagram

| No. | Revision/Issue | Date |
|-----|----------------|------|
| | | |



200 Broadhollow Road-Suite 207
Melville, New York 11747

Project Name and Address

840 East 233rd Street
Bronx, NY 10466

Project 9628-BXNY

Date 3/11/2025

Scale As Noted

Figure

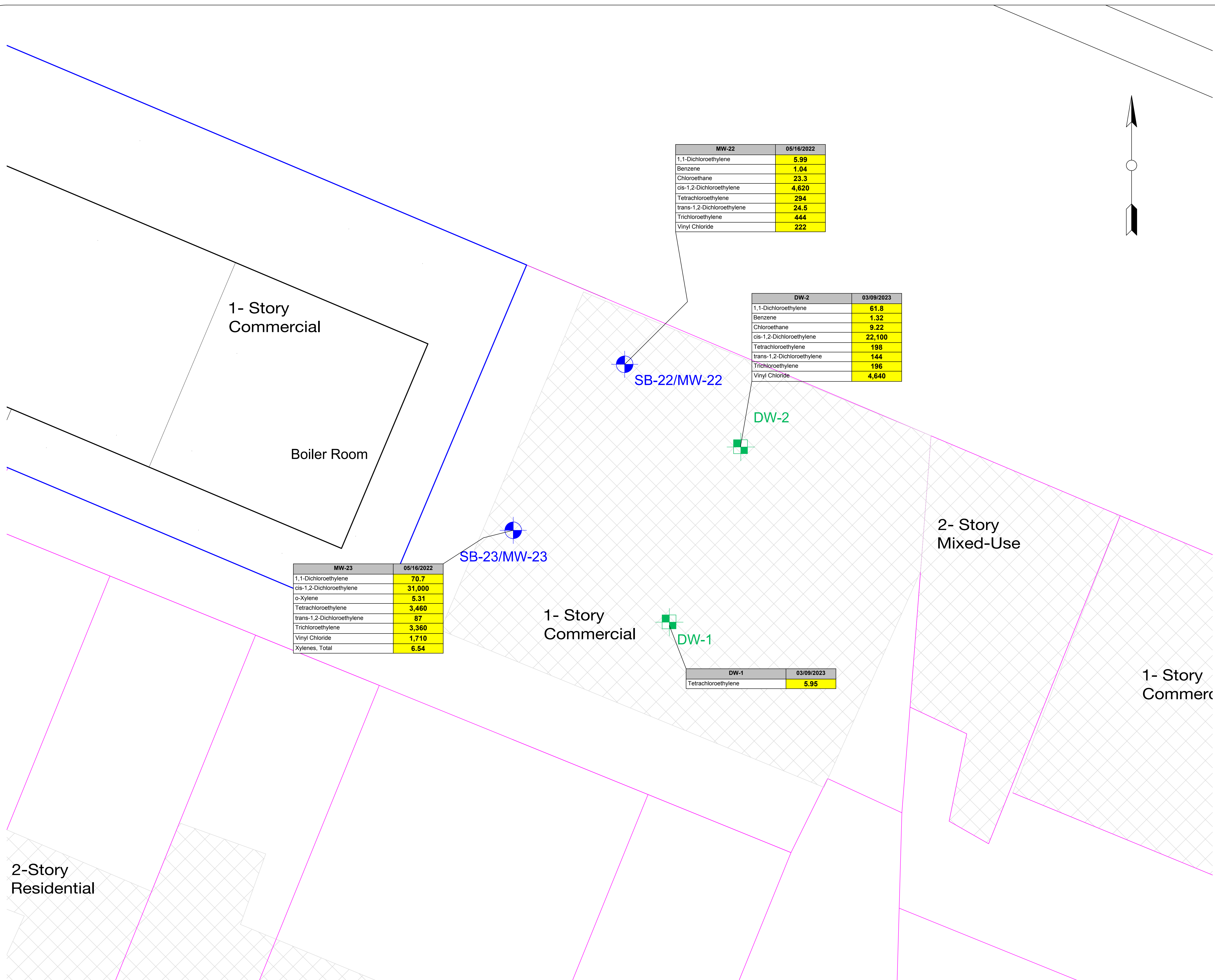
10

| MW-22 | 05/16/2022 |
|----------------------------|------------|
| 1,1-Dichloroethylene | 5.99 |
| Benzene | 1.04 |
| Chloroethane | 23.3 |
| cis-1,2-Dichloroethylene | 4,620 |
| Tetrachloroethylene | 294 |
| trans-1,2-Dichloroethylene | 24.5 |
| Trichloroethylene | 444 |
| Vinyl Chloride | 222 |

| DW-2 | 03/09/2023 |
|----------------------------|------------|
| 1,1-Dichloroethylene | 61.8 |
| Benzene | 1.32 |
| Chloroethane | 9.22 |
| cis-1,2-Dichloroethylene | 22,100 |
| Tetrachloroethylene | 198 |
| trans-1,2-Dichloroethylene | 144 |
| Trichloroethylene | 196 |
| Vinyl Chloride | 4,640 |

| MW-23 | 05/16/2022 |
|----------------------------|------------|
| 1,1-Dichloroethylene | 70.7 |
| cis-1,2-Dichloroethylene | 31,000 |
| o-Xylene | 5.31 |
| Tetrachloroethylene | 3,460 |
| trans-1,2-Dichloroethylene | 87 |
| Trichloroethylene | 3,360 |
| Vinyl Chloride | 1,710 |
| Xylenes, Total | 6.54 |

| DW-1 | 03/09/2023 |
|---------------------|------------|
| Tetrachloroethylene | 5.95 |



**Appendix A:
IRMWP Approval**

NEW YORK STATE DEPARTMENT OF ENVIRONMENTAL CONSERVATION

Division of Environmental Remediation, Region 2
47-40 21st Street, Long Island City, NY 11101
P: (718) 482-4995
www.dec.ny.gov

April 22, 2022

Suhak Lee
Attn: Jacob B. Lee
150 County Road, #29
Tenafly, NJ 07670

**Re: Jay Dry Cleaners
840 East 233rd Street, Bronx, Bronx County
Brownfield Cleanup Program # C203119
Interim Remedial Measure Work Plan**

Dear Mr. Lee:

The New York State Department of Environmental Conservation (NYSDEC), in consultation with the New York State Department of Health, has completed its review of the Interim Remedial Measure Work Plan (IRMWP) dated February 4, 2022, which was prepared by P.W. Grosser Consulting Engineer & Hydrogeologist, P.C. on behalf of Suhak Lee (the Participant). The IRMWP was submitted to NYSDEC under the Brownfield Cleanup Program. The IRMWP is deemed adequate and is hereby approved for implementation.

The Participant and its contractors are solely responsible for safe execution of all invasive and other field work performed under this IRMWP. The Participant and its contractors must obtain all local, state, and/or federal permits or approvals that may be required to perform work under this IRMWP. Further, the Participant and its contractors are solely responsible for the identification of utilities that might be affected by work under this IRMWP and, implementation of all required, appropriate, or necessary health and safety measures during performance of work under this IRMWP.

In accordance with the requirements of the Brownfield Cleanup Agreement and the Citizen Participation Plan, the approved IRMWP must be placed in the project document repositories within 5 business days. Any draft copies of this work plan should be removed. A certification that this document has been placed in the project repositories, and that the repositories are complete with all project documents, must be submitted to the NYSDEC project manager.

Please notify NYSDEC at least 7 days prior to commencing any field work related to the approved IRMWP. Should you have any questions regarding this letter or any other aspect of the project, please contact me at 718-482-7541 or wendi.zheng@dec.ny.gov.

Regards,

Wendi Zheng

Wendi Zheng
Project Manager

cc: J. O'Connell, C. Maycock, H. Leibowitz – NYSDEC
S. McLaughlin, E. O'Neil – NYSDOH
P. Boyce, R. Morley – PWGC
P. Stewart, J. Stewart – Advanced Cleanup Technologies, Inc.

Appendix B: Photo Log



Existing exposed trenches to be sealed during IRM



CAMP Equipment



Dri-Eaz HEPA 700 Air Scrubber with an activated carbon filter



Ventilation Fan during ground intrusive work



Bosch HEPA Dust Collector to collect dust during concrete trench work



Expanded Trenches with Sealed drains and subsurface piping to be installed



Expanded Trenches with Sealed drains and subsurface piping to be installed



Existing concrete gouges that were sealed



Installed Catch basins to replaced brick exposed sumps



Containerized concrete and soil waste



Newly installed catch basin following installation



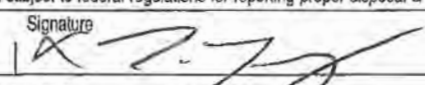
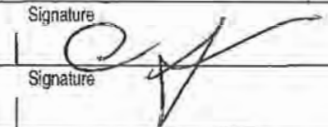
Newly installed drains following installation



Dewatering Well with Pump

**Appendix C:
Waste Manifests**

Please print or type
(Form designed for use on olive (12-pitch) typewriter.)

| | | | | | | | |
|--|---|--|--|--|--------------------|-------------------|------|
| NON-HAZARDOUS WASTE MANIFEST | 1. Generator ID Number N Y S C E S G G | 2. Page 1 of 1 | 3. Emergency Response Phone 5167813000 | 4. Waste Tracking Number 2023-8258-ACT | | | |
| 5. Generator's Name and Mailing Address COMMERCIAL PROPERTY 840 E 233RD ST BRONX NY 10468 | | Generator's Site Address (if different than mailing address) 864 E 233RD ST BRONX NY 10468 | | | | | |
| Generator's Phone: 516 840-2947 | | | | | | | |
| 6. Transporter 1 Company Name Action Trucking Inc. | | U.S. EPA ID Number NYD064748304 | | | | | |
| 7. Transporter 2 Company Name Veolia ES Technical Solutions | | U.S. EPA ID Number NJD060831380 | | | | | |
| 8. Designated Facility Name and Site Address Veolia ES Technical Solutions Inc 1 Eden Lane Flanders NJ 07838 | | U.S. EPA ID Number NJD080538593 | | | | | |
| Facility's Phone: 973 347-1309 | | | | | | | |
| GENERATOR | 9. Waste Shipping Name and Description | | 10. Containers | | 11. Total Quantity | 12. Unit Wt./Vol. | |
| | | | No. | Type | | | |
| | 1. Non-RCRA, non-DOT Regulated Material (Soil Borings and Drilling Spoils) | | 003 | DM | 1725 | F | |
| | 2. | | | | | | |
| | 3. | | | | | | |
| 4. | | | | | | | |
| 13. Special Handling Instructions and Additional Information (1) WIP # 516 (ID 27, L) (ADVANCED CLEANUP TECHNOLOGIES) | | | | | | | |
| 14. GENERATOR'S CERTIFICATION: I certify the materials described above on this manifest are not subject to federal regulations for reporting proper disposal of Hazardous Waste. | | | | | | | |
| Generator's/Officer's Printed/Typed Name Tommy Act | | Signature  | | Month 13 | Day 1 | Year 23 | |
| TRANSPORTER | 15. International Shipments <input type="checkbox"/> Import to U.S. <input type="checkbox"/> Export from U.S. | | Port of entry/exit Date leaving U.S.: | | | | |
| | Transporter Signature (for exports only): | | | | | | |
| 16. Transporter Acknowledgment of Receipt of Materials | | | | | | | |
| Transporter 1 Printed/Typed Name Charles Fowser | | Signature  | | Month 13 | Day 1 | Year 23 | |
| Transporter 2 Printed/Typed Name | | Signature | | Month | Day | Year | |
| DESIGNATED FACILITY | 17. Discrepancy | | | | | | |
| | 17a. Discrepancy Indication Space <input type="checkbox"/> Quantity <input type="checkbox"/> Type <input type="checkbox"/> Residue <input type="checkbox"/> Partial Rejection <input type="checkbox"/> Full Rejection | | | | | | |
| | 17b. Alternate Facility (or Generator) | | | Manifest Reference Number: | | | |
| | Facility's Phone: | | | U.S. EPA ID Number | | | |
| 17c. Signature of Alternate Facility (or Generator) | | | | | Month | Day | Year |
| 18. Designated Facility Owner or Operator: Certification of receipt of materials covered by the manifest except as noted in Item 17a | | | | | | | |
| Printed/Typed Name | | Signature | | Month | Day | Year | |

Please print or type
 (Form available for use on site (12-0105) typewriter.)

| | | | | | |
|--|--|--|--|--|--|
| NON-HAZARDOUS WASTE MANIFEST | | 1. Generator ID Number CESQG | 2. Page 1 of 1 | 3. Emergency Response Phone 516-781-3000 | 4. Waste Tracking Number 2024-1337-ACT |
| 5. Generator's Name and Mailing Address Commercial Property 840 E 233rd Street Bronx, NY 10466 Generator's Phone: 516-640-2947 | | | Generator's Site Address (if different than mailing address) 864 E. 233 RD. Bronx NY 10466 | | |
| 6. Transporter 1 Company Name Action Trucking Co., Inc. | | Phil#: 516-781-3000 State ID#: - | | U.S. EPA ID Number NYD064749304 | |
| 7. Transporter 2 Company Name Veolia ES Technical Solutions | | Phil#: 973-347-7111 | | U.S. EPA ID Number NJD080631369 | |
| 8. Designated Facility Name and Site Address Veolia ES Technical Solutions Inc 1 Eden Lane Flanders, NJ 07838 Facility's Phone: 973-347-1909 | | | State ID#: NJ- | | U.S. EPA ID Number NJD980536593 |

| 9. Waste Shipping Name and Description | 10. Containers | | 11. Total Quantity | 12. Unit Wt./Vol. | L | ID27 |
|---|----------------|-----------|--------------------|-------------------|---|------|
| | No. | Type | | | | |
| 1. NONRCRA / NONDOT REGULATED MATERIAL (SOIL BORINGS AND DRILLING SPOILS) <i>SOLID</i> | 001 | DM | 625 | P | | |
| 2. | | | | | | |
| 3. | | | | | | |
| 4. | | | | | | |

13. Special Handling Instructions and Additional Information
01: () Soil Borings and Drilling Spoils *55*
WDP# 11439002
(Advanced Technologies)
Delay received 6/18/24

14. GENERATOR'S CERTIFICATION: I certify the materials described above on this manifest are not subject to federal regulations for reporting proper disposal of Hazardous Waste.
 Generator's/Officer's Printed/Typed Name: *Christopher Carr* Signature: *[Signature]* Month: **6** Day: **11** Year: **24**

15. International Shipments Import to U.S. Export from U.S. Port of entry/exit: _____ Date leaving U.S.: _____

16. Transporter Acknowledgment of Receipt of Materials
 Transporter 1 Printed/Typed Name: *Carlos Fonseca* Signature: *[Signature]* Month: **6** Day: **11** Year: **24**
 Transporter 2 Printed/Typed Name: *Joshua Stewart* Signature: *[Signature]* Month: **6** Day: **17** Year: **24**

17. Discrepancy
 17a. Discrepancy Indication Space: Quantity Type Residue Partial Rejection Full Rejection
 Manifest Reference Number: _____

17b. Alternate Facility (or Generator) U.S. EPA ID Number: _____
 Facility's Phone: _____
 17c. Signature of Alternate Facility (or Generator) Month: _____ Day: _____ Year: _____

18. Designated Facility Owner or Operator: Certification of receipt of materials covered by the manifest except as noted in Item 17a
 Printed/Typed Name: *DEANNA PARLUE* Signature: *[Signature]* Month: **6** Day: **18** Year: **24**

**Appendix D:
CAMP and Field Notes**

DirtTrak
rental

Trenching Day 1
Drains
corrugated

ADVANCED CLEANUP TECHNOLOGIES
COMMUNITY AIR MONITORING LOG

8/22/22

Project/Address: 9628-BXNY dft. re Basecoat

Weather: (AM) rain humid Temperature: (AM) 70's Wind Direction (AM) _____
(PM) cloudy humid (PM) 80's (PM) _____

| <u>Location</u> | <u>Time</u> | <u>PID Reading</u> | <u>PID</u> |
|-----------------|-------------|----------------------------|------------|
| Basecoat | 11:00 AM | PDR | PID |
| pre-trench | ↳ | 0.058 mg/m ³ | Oppb |
| trenching | 10:30 | 0.058 mg/m ³ | Oppb |
| break | 11:30 | 0.049 mg/m ³ | Oppb |
| trenching | 12:00 | 0.357 mg/m ³ | Oppb |
| " | 12:30 | 0.431 mg/m ³ | Oppb |
| " | 1 pm | 0.065 mg/m ³ | Oppb |
| " | 2 pm | 0.580 mg/m ³ | Oppb |
| break | 2:45 | .060 | Oppb |
| pack up | 3 pm | .050 | Oppb |
| End | 3:30 | .056 | Oppb |

Monitor: Tim Yang
Name

[Signature]
Signature

Vent fan + HEPA scrubber
Running

PDR
PDR meter
DustTrak

PID ACT
ppb

HEPA scrubber
running

ADVANCED CLEANUP TECHNOLOGIES
COMMUNITY AIR MONITORING LOG

8/23/22

Project/Address: 9628-BXNY offsite

Weather: (AM) hot Temperature: (AM) 80 Wind Direction (AM) basement
(PM) _____ (PM) _____ (PM) _____

| <u>Location</u> Activity | <u>Time</u> | <u>PID Reading</u> | |
|------------------------------------|-------------|---------------------------|------|
| | | PDR | PID |
| mobile 9-10 | | | |
| Trenching | 1000 | .208 mg/m ³ | 0ppb |
| " | 1100 | .112 mg/m ³ | 0ppb |
| " | 1200 | .010 mg/m ³ | 0ppb |
| break | 1300 | .222 mg/m ³ | 0ppb |
| Trenching PDR batteries dead | 1400 | - | 0ppb |
| attach drain fittings | 1500 | - | 0ppb |
| | | | |
| | | | |
| | | | |
| | | | |
| | | | |
| | | | |
| | | | |
| | | | |

Monitor: Tim Yung
Name

[Signature]
Signature

HEPA scrubber
running

ADVANCED CLEANUP TECHNOLOGIES
COMMUNITY AIR MONITORING LOG

8/24/22

Project/Address: 9628 - BXNY

Weather: (AM) hot Temperature: (AM) 80's Wind Direction (AM) _____
(PM) _____ (PM) _____ (PM) _____

Location Time PID Reading

Only @ site a few hours
no trenching or concrete breaking
attempts to seal fittings on
Drains pieces
fill bathroom drain w/ gravel
fit drain on end

Didn't set up CAMP- fieldwork cut short

TY

Monitor: TY
Name

Signature

ADVANCED CLEANUP TECHNOLOGIES
COMMUNITY AIR MONITORING LOG

9/7/22

Project/Address: 9628-BXNY off site

Weather: (AM) cool rainy Temperature: (AM) ~70's Wind Direction (AM) basement
(PM) _____ (PM) _____ (PM) _____

| Location Activity | Time | PID Reading | |
|----------------------|------|------------------------------|--------|
| | | PDR $\mu\text{g}/\text{m}^3$ | PID |
| mobilize | 0915 | conc 0.013 TWA 0.010 | 0 ppb |
| Filling Tanks | 1000 | history dead ✓ replace | 0 ppb |
| prep for trench fill | 1015 | conc .025 TWA .031 | 0 ppb |
| " " | 1100 | conc .000 TWA .006 | 0 ppb |
| same | 1130 | conc .006 TWA .001 | 0 ppb |
| concret | 1245 | conc .220 TWA .095 | 70 ppb |
| concrete | 1330 | conc .142 TWA .121 | 80 ppb |
| pack up finish up | 1405 | conc .144 TWA .116 | 95 ppb |

Monitor: Tim Young ACT
Name

[Signature]
Signature

PDR - pine
thermo
scientific
PDR

PID - ACT
PDR
Rae

HEPA air
scrubber running
entire time

9/8/22

ADVANCED CLEANUP TECHNOLOGIES
COMMUNITY AIR MONITORING LOG

Project/Address: 9628-BANY cfs, k

Weather: (AM) _____ Temperature: (AM) _____ Wind Direction (AM) _____
(PM) _____ (PM) _____ (PM) _____

| Location | Time | PDR | PID Reading | PID |
|-----------------|------|------|------------------------|--------|
| Swamp walk | 1000 | 1 | - | |
| Swamp + grant | 1100 | .081 | | 48 ppb |
| | | .091 | | |
| " | 1200 | .031 | | 63 ppb |
| | | .100 | | |
| " | 1240 | .232 | | 53 ppb |
| | | .137 | | |
| " | 1330 | .091 | | 60 ppb |
| catch basin | | .133 | | |
| trash (pile) | 1400 | .328 | | 54 ppb |
| | | .179 | | |
| " | 1430 | .099 | | 50 ppb |

* PID needs zero/ calibration

Monitor: Tim Yung ACT
Name

[Signature]
Signature

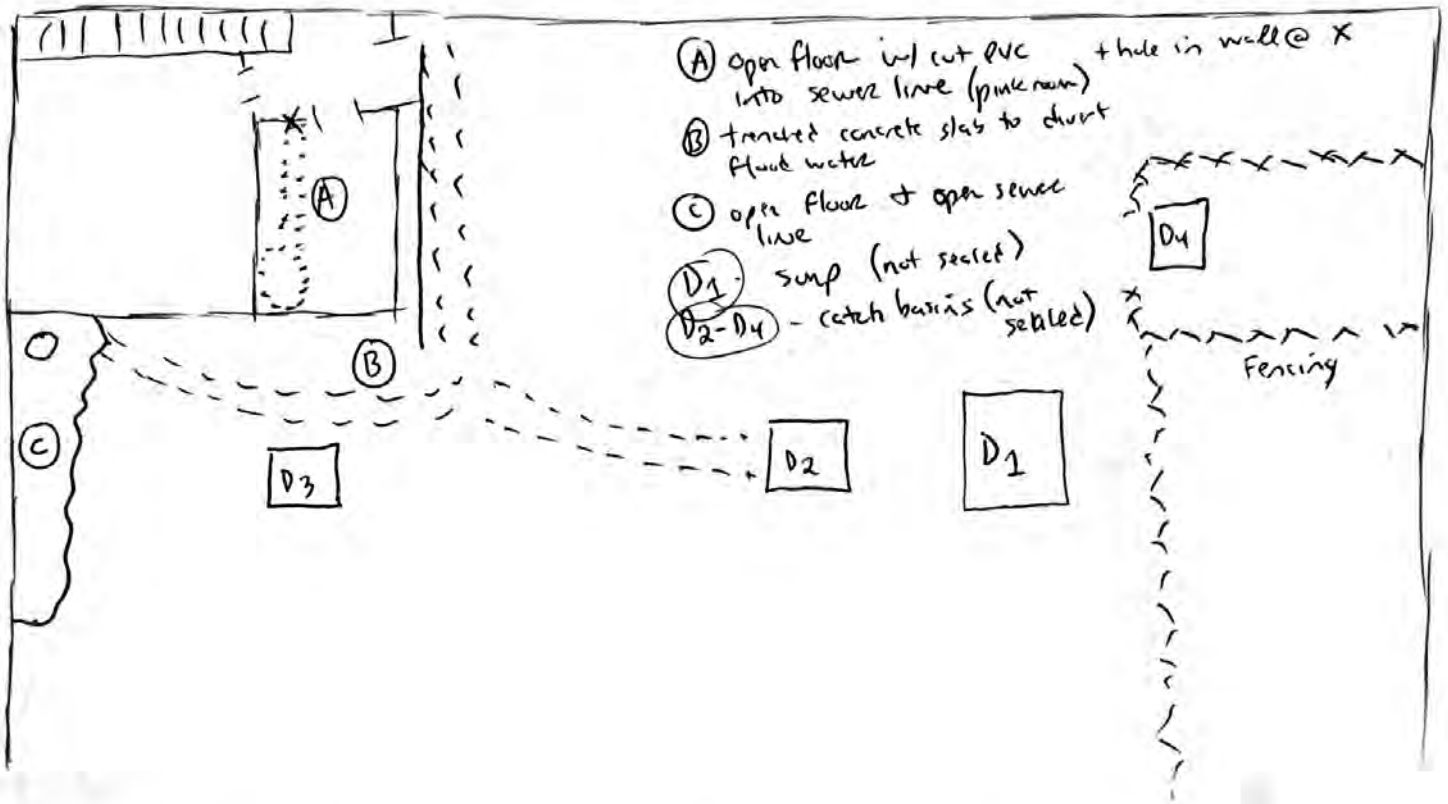
PDR - Pine Thermo Scientific
PDR

PID - Act Pph AAE

HEPA air scrubber running entire time

9628-BXNY
 offsite - drainage - remediation
Before Drains, sump, sealing floor

(TY)



After ACT drains, sewer drains, sump, sealing floor

- (A) Previous PVC swapped for new 3" horizontal w/ Drain on 45° @ X
 Drain sealed w/ hydraulic cement + non shrink grout
 PVC backfilled w/ gravel - sealed w/ high strength
 PVC connected to previous open in sewer line
 Tested ✓
- (B) open floor trenching replaced w/ channel drains - set in
 trench on bedding mason sand, backfilled gravel + sealed w/
 non shrink grout - connected to each section, T + catch
 basins w/ sealed fittings + 4" corrugated piping
- (C) Drain installed (as depression) over open sewer line - open ~~floor~~
 broken concrete, chipped out, cleaned up - bedded w/ grout
 molded on slopes to depression drain w/ concrete
 tested ✓

D1
 open sump replaced w/ sealed sump
 set in, backfilled, sealed w/ grout
 pipe tubes from D2 + D4 - sump re connected
 tested ✓

D2 + D4 - open catch basins
 replaced w/ sealed catch basins
 connecting to new sump

D3 - open catch basin backfilled
 + sealed w/ grout

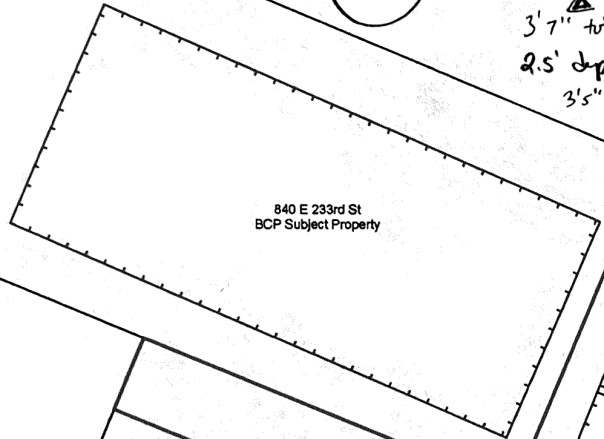


TY

3/1/23 - Drum pickup by Action
 3 drums
 install DW-1
 3' 7" total depth
 2.5' depth water
 3.5" 4" well screen
 installed
 low PID

3/2/23
 install DW-2
 4' 2" total depth
 to BR (assumed)
 3' depth to water
 4" 4" well screen
 installed

high PIDs - 14 @ surface
 (in hole)
 20ppm @ 3'
 32ppm @ 4'
 0-100pph outside the hole



Proposed Post IRM Air Sample

- Proposed Post IRM Air Sample
- Trench
- drains to sump
- Interior Wall
- Building Outline
- Tax Lot
- Site Boundary

PWGC
 CLIENT DRIVEN SOLUTIONS
 P. W. Grosser Consulting, Inc.
 630 Johnson Ave., Suite 7
 Bohemia, NY 11716
 Ph: 631-589-8353 • Fax: 631-589-8705
 pwgc.info@pwgrosser.com

UNAUTHORIZED REVISION OR ADDITION TO THE DRAWING AND RELATED DOCUMENTS IS A VIOLATION OF SEC. 1709 OF THE E.P.C.L. OR OTHER LAW.
 DRAWING PREPARED FOR:

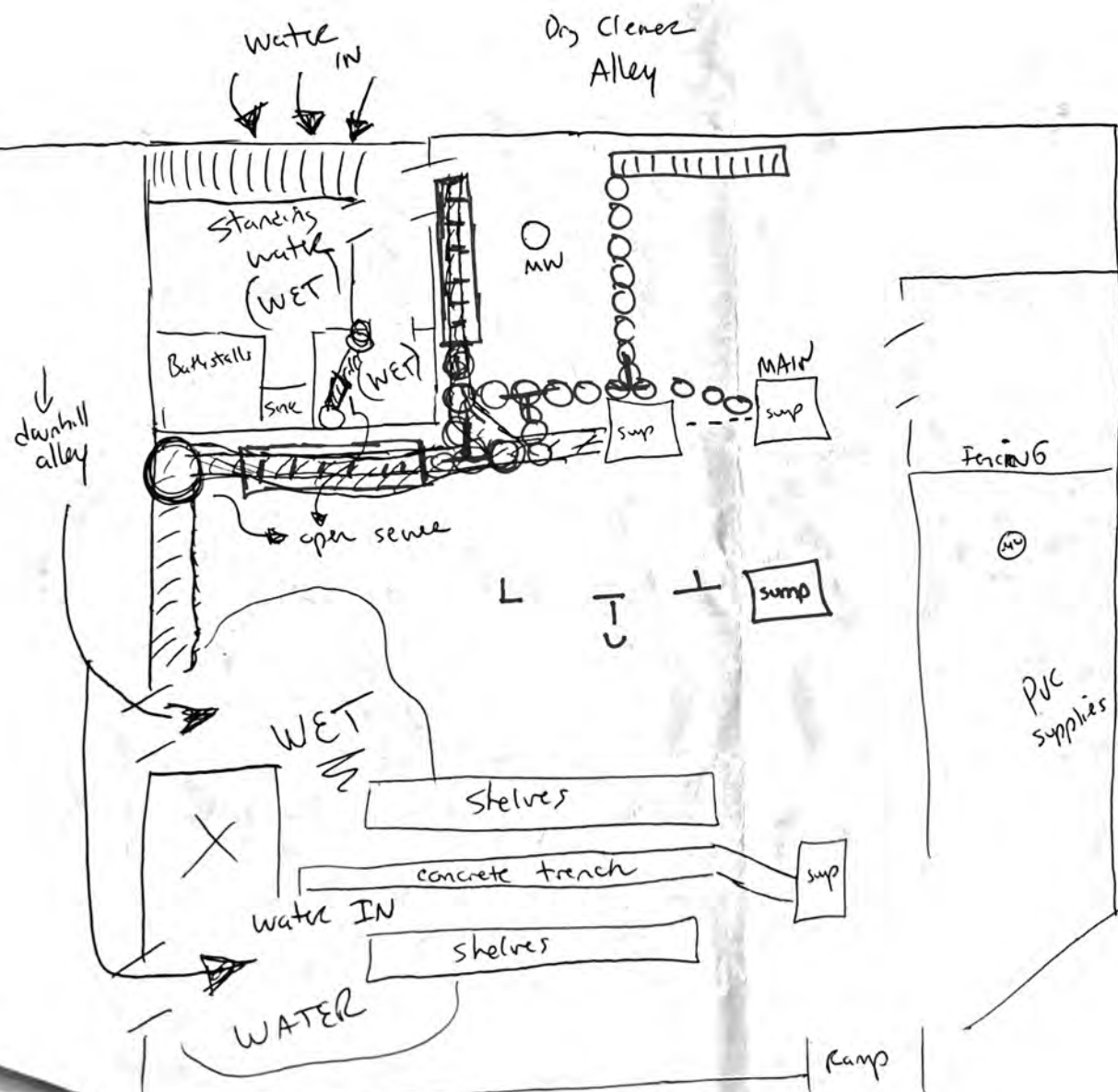
| REVISION | DATE | INITIAL | COMMENTS |
|----------|------|---------|----------|
| | | | |
| | | | |
| | | | |
| | | | |
| | | | |
| | | | |
| | | | |
| | | | |
| | | | |
| | | | |

DRAWING INFORMATION:
 Project: ACT2101 Designed by: BM
 Date: 12/15/2022 Drawn by: PH
 Scale: AS SHOWN Approved by: BM

SITE PLAN

866 E 233rd St.
 Bronx, NY

FIGURE NO:
 2



10' drain section (L)

4" corrugated

adapted T's

sump

110 Main Street, Port Washington, NY

Date: 8/9/23

SAMPLE INFORMATION RECORD

PROJECT NO.: 9628-BXNY Sampling Personnel: TY
Job Locations: Offsite Sampling (Chris. D'Amico) 233rd Barry

Field Sample Designation: IA-3 Time: _____

SAMPLE TYPE:

GROUNDWATER: _____ SEDIMENT: _____

SURFACE WATER: _____ SOIL: _____

AIR (specify): _____ (soil Vapor, Indoor air, Outdoor air) OTHER (describe): _____

GROUNDWATER INFORMATION:

Depth to Groundwater: _____ Measurement Method: _____

Depth of well or Sampling Point: _____ Measurement Method: _____

Volume of Groundwater Purged: _____ Purge Method: no PID available

FIELD TEST RESULTS:

Color: _____ pH: _____ Odor: _____

Temperature (°F/°C): _____ Specific Conductance (umhos/cm): _____

Other: PID: _____ ppm Helium Detector: _____ ppm

Canister #: 23348 Initial Pressure: -36 8/9 Start Time: 1430 8/9

Flow Regulator #: 2860 Final Pressure: ~~0~~ 8/10 End Time: 0910 8/10

- 9

SAMPLE ANALYSIS:

REMARKS: IA - 24 HR closest to dry cleaned
by the air scrubber (currently off)
removed

Date: 8/9/23

SAMPLE INFORMATION RECORD

PROJECT NO.: 9628-BXNY Sampling Personnel: TY

Job Locations: Offsite Sampling

Field Sample Designation: IA-4 Time: _____

SAMPLE TYPE:

GROUNDWATER: _____

SEDIMENT: _____

SURFACE WATER: _____

SOIL: _____

AIR (specify): _____ (soil Vapor, Indoor air, Outdoor air)

OTHER (describe): _____

GROUNDWATER INFORMATION:

Depth to Groundwater: _____

Measurement Method: _____

Depth of well or Sampling Point: _____

Measurement Method: _____

Volume of Groundwater Purged: _____

Purge Method: no PID available

FIELD TEST RESULTS:

Color: _____ pH: _____ Odor: _____

Temperature (°F/°C): _____ Specific Conductance (µmhos/cm): _____

Other: PID: _____ ppm Helium Detector: _____ ppm

Canister #: 28570 Initial Pressure: -30 8/9 Start Time: 1425 8/9

Flow Regulator #: 2854 Final Pressure: -10 8/10 End Time: 0945 8/10

SAMPLE ANALYSIS:

REMARKS:

IA - 24 hr - mid. basement

110 Main Street, Port Washington, NY

Date: 8/9/23

SAMPLE INFORMATION RECORD

PROJECT NO.: 9628-BXNY Sampling Personnel: TY

Job Locations: Offsite Sampling E23312 Bronx

Field Sample Designation: IA-7 Time: _____

SAMPLE TYPE:

GROUNDWATER: _____ SEDIMENT: _____

SURFACE WATER: _____ SOIL: _____

AIR (specify): _____ (soil Vapor, Indoor air, Outdoor air) OTHER (describe): _____

GROUNDWATER INFORMATION:

Depth to Groundwater: _____ Measurement Method: _____

Depth of well or Sampling Point: _____ Measurement Method: _____

Volume of Groundwater Purged: _____ Purge Method: no PID available

FIELD TEST RESULTS:

Color: _____ pH: _____ Odor: _____

Temperature (°F/°C): _____ Specific Conductance (µmhos/cm): _____

Other: PID: _____ ppm Helium Detector: _____ ppm

Canister #: 489 Initial Pressure: -30 8/9/23 Start Time: 1420 8/9/23

Flow Regulator #: 2988 Final Pressure: -7 8/10/23 End Time: 0900 8/10/23

SAMPLE ANALYSIS:

REMARKS: IA - 24 hr - farthest East from dry clean



587 East Middle Turnpike, P.O. Box 370, Manchester, CT 06040
Telephone: 860.645.1102 • Fax: 860.645.0623

CHAIN OF CUSTODY RECORD

AIR ANALYSES

860-645-1102

email: greg@phoenixlabs.com

P.O. #

Page 1 of 1

Data Delivery:

Fax #: _____
 Email: _____
 Phone #: _____

ACT

| | | |
|--|---|--|
| Report to: Jason Stewart | Project Name: 9268-BXNY | Data Format: (Circle) Equis Excel Other: _____ |
| Customer: Advanced Cleanup Technologies, Inc. | Invoice to: Karen Friedman | Requested Deliverable: RCP ASP CAT B |
| Address: 228 Park Ave S PMB 34864 | | MCP NJ Deliverables |
| 13763 New York, New York 10003 | Sampled by: Yisung Yang / Ray Buzeta | Quote Number: _____ |

| Phoenix ID # | Client Sample ID | Canister ID # | Canister Size (L) | Outgoing Canister Pressure ("Hg) | Incoming Canister Pressure ("Hg) | Flow Regulator ID # | Flow Controller Setting (mL/min) | Sampling Start Time | Sampling End Time | Sample Start Date | Canister Pressure at Start ("Hg) | Canister Pressure at End ("Hg) | Ambient/Indoor Air | Soil Gas | Grab (G) Composite (C) | TO-15 | APH |
|--------------------------------------|------------------|---------------|-------------------|----------------------------------|----------------------------------|---------------------|----------------------------------|---------------------|-------------------|-------------------|----------------------------------|--------------------------------|--------------------|----------|------------------------|-------|-----|
| THIS SECTION FOR LAB USE ONLY | | | | | | | | | | | | | | | | | |
| 92021 | IA-4 | 28560 | 6.0L | -30 | 40 | 5043 | 3.6 | 15:00 8/1/22 | 11:05 8/2/22 | 8/1/22 | -30 | -8 | ✓ | | | ✓ | |
| | | 23338 | 6.0L | -30 | | 5042 | 3.6 | | | | | | | | | | |
| 92022 | IA-7 | 23330 | 6.0L | -30 | 40 5 | 5000 | 3.8 | 15:00 8/1/22 | 10:22 8/2/22 | 8/1/22 | -30 | -6 | ✓ | | | ✓ | |

| | | | | |
|--|---------------------------------|---------------------|--------------------|---|
| Relinquished by: AET Ray Buzeta | Accepted by: [Signature] | Date: 8-3-22 | Time: 10:54 | I attest that all media released by Phoenix Environmental Laboratories, Inc. have been received in good working condition and agree to the terms and conditions as listed on the back of this document. |
| | | Date: 8/3/22 | Time: 1750 | |

| | | | | | | | |
|---|--|--|-------------------------|-------------------------|-----------------|-------------------------|-------------------------|
| State Where Samples Collected: _____ | Turnaround Time: 1 Day <input type="checkbox"/> | Requested Criteria: (Please Circle) CT: _____ | MA: _____ | NI: _____ | NY: _____ | PA: _____ | VT: _____ |
| SPECIAL INSTRUCTIONS, OC REQUIREMENTS, REGULATORY INFORMATION: Can # 23338 (3) - 6.0L 24 hr returned unused | 2 Day <input type="checkbox"/> | TAC I/C | Indoor Air: Residential | Indoor Air: Residential | Vapor Intrusion | Indoor Air: Residential | Indoor Air: Residential |
| | 3 Day <input type="checkbox"/> | TAC RES | Ind/Commercial | Ind/Commercial | | Residential | Residential |
| | 4 Day <input type="checkbox"/> | SVVC I/C | Soil Gas: Residential | Soil Gas: Residential | | Non-residential | Industrial Sub-slab |
| | 5 Day <input type="checkbox"/> | SVVC RES | Residential | Residential | | | Residential |
| | | | GWV I/C | Ind/Commercial | Ind/Commercial | | |

Data Delivery:

Fax #:

Email:

Phone #:

| | | |
|--|-----------------------------------|---|
| Report to: Ray Buzeta | Project Name: 9628-BXNY | Data Format: (Circle) Equis Excel Other: |
| Customer: Advanced Cleanup Technologies, Inc. | Invoice to: Karen Friedman | Requested Deliverable: RCP ASP CAT B |
| Address: 228 Park Ave S FMB 34864 | | MCP NJ Deliverables |
| 14905 New York, New York 10003 | Sampled by: Tim Young | Quote Number: |

| Phoenix ID # | Client Sample ID | Canister ID # | Canister Size (L) | Outgoing Canister Pressure ("Hg) | Incoming Canister Pressure ("Hg) | Flow Regulator ID # | Flow Controller Setting (mL/min) | Sampling Start Time | Sampling End Time | Sample Start Date | Canister Pressure at Start ("Hg) | Canister Pressure at End ("Hg) | Ambient/Indoor Air | Soil Gas | Grab (G) Composite (C) | TO-15 | APH |
|-------------------------------|------------------|---------------|-------------------|----------------------------------|----------------------------------|---------------------|----------------------------------|---------------------|-------------------|-------------------|----------------------------------|--------------------------------|-------------------------------------|----------|------------------------|-------------------------------------|-----|
| THIS SECTION FOR LAB USE ONLY | | | | | | | | | | | | | | | | | |
| | IA-4 | 28570 | 6.0L | -30 | | 2854 | 3.01 | 1425 | 0945 | 8/6/23 | -30 | -10 | <input checked="" type="checkbox"/> | | | <input checked="" type="checkbox"/> | |
| | IA-3 | 23348 | 6.0L | -30 | | 2860 | 3.17 | 1430 | 0910 | 8/6/23 | -30 | -9 | <input checked="" type="checkbox"/> | | | <input checked="" type="checkbox"/> | |
| | IA-7 | 489 | 6.0L | -30 | | 2988 | 3.28 | 1420 | 0900 | 8/6/23 | -30 | -7 | <input checked="" type="checkbox"/> | | | <input checked="" type="checkbox"/> | |
| | | | | | | | | ↑ | ↑ | | | | | | | | |
| | | | | | | | | 8/6/23 | 8/10/23 | | | | | | | | |

| | | | | |
|------------------|--------------|-------|-------|---|
| Relinquished by: | Accepted by: | Date: | Time: | I attest that all media released by Phoenix Environmental Laboratories, Inc. have been received in good working condition and agree to the terms and conditions as listed on the back of this document: |
| | | | | Signature: [Signature] Date: 8/10/23 |

| | | | | | | |
|--|---|---|---|-------------------------------------|--|--|
| State Where Samples Collected: Brnx, NY | Turnaround Time: 1 Day* <input type="checkbox"/> 2 Day* <input type="checkbox"/> 3 Day* <input type="checkbox"/> 4 Day* <input type="checkbox"/> 5 Day* <input type="checkbox"/> Standard <input checked="" type="checkbox"/> | Requested Criteria: TAC I/C TAC RES SVVC I/C SVVC RES GWV I/C GWV CES | Please Circle: MA: Indoor Air Residential Ind/Commercial Soil Gas: Residential Ind/Commercial | NI: NY Vapor Intrusion | PA: Indoor Air Residential Non-residential | VT: Indoor Air Residential Industrial Sub-slab Residential Industrial |
| SPECIAL INSTRUCTIONS, QC REQUIREMENTS, REGULATORY INFORMATION: (3) - 6.0L 24 hr | *SURCHARGES MAY APPLY | | | | | |

Appendix E: Laboratory Reports



Wednesday, January 10, 2024

Attn:
Advanced Cleanup Technologies, Inc.
228 Park Ave S PMB 34864
New York, New York 10003

Project ID: 9268-BXNY
SDG ID: GCL97021
Sample ID#s: CL97021 - CL97022

This laboratory is in compliance with the NELAC requirements of procedures used except where indicated.

This report contains results for the parameters tested, under the sampling conditions described on the Chain Of Custody, as received by the laboratory. This report is incomplete unless all pages indicated in the pagination at the bottom of the page are included.

A scanned version of the COC form accompanies the analytical report and is an exact duplicate of the original.

Enclosed are revised Analysis Report pages. Please replace and discard the original pages. If you are the client above and have any questions concerning this testing, please do not hesitate to contact Phoenix Client Services at ext.200. The contents of this report cannot be discussed with anyone other than the client listed above without their written consent.

Sincerely yours,

A handwritten signature in black ink that reads "Phyllis Shiller". The signature is written in a cursive style with a large initial "P".

Phyllis Shiller
Laboratory Director

NELAC - #NY11301
CT Lab Registration #PH-0618
MA Lab Registration #M-CT007
ME Lab Registration #CT-007
NH Lab Registration #213693-A,B

NJ Lab Registration #CT-003
NY Lab Registration #11301
PA Lab Registration #68-03530
RI Lab Registration #63
VT Lab Registration #VT11301



Environmental Laboratories, Inc.
587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045
Tel. (860) 645-1102 Fax (860) 645-0823



**NY ANALYTICAL SERVICES PROTOCOL
DATA PACKAGE**

Client: Advanced Cleanup Technologies, Inc.

Project: 9268-BXNY

Laboratory Project: GCL97021



Environmental Laboratories, Inc.
587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06040
Tel. (860) 645-1102 Fax (860) 645-0823



NY Analytical Services Protocol Format

January 10, 2024

SDG I.D.: GCL97021

Advanced Cleanup Technologies, Inc. 9268-BXNY

Methodology Summary

Volatiles in Air

Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air: Method TO-15, Second Edition, U. S. Environmental Protection Agency, January 1999.



Environmental Laboratories, Inc.
587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06040
Tel. (860) 645-1102 Fax (860) 645-0823



NY Analytical Services Protocol Format

January 10, 2024

SDG I.D.: GCL97021

Advanced Cleanup Technologies, Inc. 9268-BXNY

Laboratory Chronicle

| Sample | Analysis | Collection Date | Prep Date | Analysis Date | Analyst | Hold Time Met |
|---------|------------------|-----------------|-----------|---------------|---------|---------------|
| CL97021 | Volatiles (TO15) | 08/02/22 | 08/03/22 | 08/03/22 | KCA | Y |
| CL97022 | Volatiles (TO15) | 08/02/22 | 08/03/22 | 08/03/22 | KCA | Y |



Environmental Laboratories, Inc.
587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045
Tel. (860) 645-1102 Fax (860) 645-0823



Sample Id Cross Reference

January 10, 2024

SDG I.D.: GCL97021

Project ID: 9268-BXNY

| Client Id | Lab Id | Matrix |
|-----------|---------|--------|
| IA-4 | CL97021 | AIR |
| IA-7 | CL97022 | AIR |



Environmental Laboratories, Inc.

587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045
Tel. (860) 645-1102 Fax (860) 645-0823



Analysis Report

January 10, 2024

FOR: Attn: Advanced Cleanup Technologies, Inc.
228 Park Ave S PMB 34864
New York, New York 10003

Sample Information

Matrix: AIR
Location Code: ACT
Rush Request: Standard
P.O.#:
Canister Id: 28560

Custody Information

Collected by: YY,RB
Received by: CP
Analyzed by: see "By" below

Date: 08/02/22 11:05
08/03/22 17:50

Laboratory Data

SDG ID: GCL97021
Phoenix ID: CL97021

Project ID: 9268-BXNY
Client ID: IA-4

Table with 11 columns: Parameter, ppbv Result, ppbv RL, LOD/MDL, ug/m3 Result, ug/m3 RL, LOD/MDL, Date/Time, By, Dilution. Rows include Volatiles (TO15) such as 1,1,1,2-Tetrachloroethane, 1,1,1-Trichloroethane, etc.

Client ID: IA-4

| Parameter | ppbv Result | ppbv RL | LOD/ MDL | ug/m3 Result | ug/m3 RL | LOD/ MDL | Date/Time | By | Dilution |
|--|----------------|------------|-------------|-----------------|-------------|-------------|-----------|-----|----------|
| Bromodichloromethane | ND | 0.149 | 0.149 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Bromoform | ND | 0.097 | 0.097 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Bromomethane | ND | 0.258 | 0.258 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Carbon Disulfide | ND | 0.321 | 0.321 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Carbon Tetrachloride | 0.085 | 0.032 | 0.032 | 0.53 | 0.20 | 0.20 | 08/03/22 | KCA | 1 |
| Chlorobenzene | ND | 0.217 | 0.217 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Chloroethane | ND | 0.379 | 0.379 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Chloroform | ND | 0.205 | 0.205 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Chloromethane | 0.596 | 0.485 | 0.485 | 1.23 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Cis-1,2-Dichloroethene | 17.2 | 0.051 | 0.051 | 68.2 | 0.20 | 0.20 | 08/03/22 | KCA | 1 |
| cis-1,3-Dichloropropene | ND | 0.221 | 0.221 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Cyclohexane | 1.81 | 0.291 | 0.291 | 6.23 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Dibromochloromethane | ND | 0.118 | 0.118 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Dichlorodifluoromethane | 0.548 | 0.202 | 0.202 | 2.71 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Ethanol | 62.8 | E 0.531 | 0.531 | 118 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Ethyl acetate | ND | 0.278 | 0.278 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Ethylbenzene | 0.928 | 0.230 | 0.230 | 4.03 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Heptane | 1.64 | 0.244 | 0.244 | 6.72 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Hexachlorobutadiene | ND | 0.094 | 0.094 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Hexane | 5.37 | 0.284 | 0.284 | 18.9 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Isopropylalcohol | 4.48 | 0.407 | 0.407 | 11.0 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Isopropylbenzene | ND | 0.204 | 0.204 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| m,p-Xylene | 3.21 | 0.230 | 0.230 | 13.9 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Methyl Ethyl Ketone | 1.70 | 0.339 | 0.339 | 5.01 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Methyl tert-butyl ether(MTBE) | ND | 0.278 | 0.278 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Methylene Chloride | ND | 0.863 | 0.863 | ND | 3.00 | 3.00 | 08/03/22 | KCA | 1 |
| n-Butylbenzene | ND | 0.182 | 0.182 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| o-Xylene | 1.23 | 0.230 | 0.230 | 5.34 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Propylene | ND | 0.581 | 0.581 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| sec-Butylbenzene | ND | 0.182 | 0.182 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Styrene | ND | 0.235 | 0.235 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Tetrachloroethene | 3.28 | 0.037 | 0.037 | 22.2 | 0.25 | 0.25 | 08/03/22 | KCA | 1 |
| Tetrahydrofuran | 6.13 | 0.339 | 0.339 | 18.1 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Toluene | 7.16 | 0.266 | 0.266 | 27.0 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Trans-1,2-Dichloroethene | ND | 0.252 | 0.252 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| trans-1,3-Dichloropropene | ND | 0.221 | 0.221 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Trichloroethene | 2.30 | 0.037 | 0.037 | 12.4 | 0.20 | 0.20 | 08/03/22 | KCA | 1 |
| Trichlorofluoromethane | 0.265 | 0.178 | 0.178 | 1.49 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Trichlorotrifluoroethane | ND | 0.131 | 0.131 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Vinyl Chloride | 1.82 | 0.078 | 0.078 | 4.65 | 0.20 | 0.20 | 08/03/22 | KCA | 1 |
| <u>QA/QC Surrogates/Internals</u> | | | | | | | | | |
| % Bromofluorobenzene | 99 | % | % | 99 | % | % | 08/03/22 | KCA | 1 |
| % IS-1,4-Difluorobenzene | 101 | % | % | 101 | % | % | 08/03/22 | KCA | 1 |
| % IS-Bromochloromethane | 100 | % | % | 100 | % | % | 08/03/22 | KCA | 1 |
| % IS-Chlorobenzene-d5 | 103 | % | % | 103 | % | % | 08/03/22 | KCA | 1 |

| Parameter | ppbv Result | ppbv RL | LOD/ MDL | ug/m3 Result | ug/m3 RL | LOD/ MDL | Date/Time | By | Dilution |
|-----------|----------------|------------|-------------|-----------------|-------------|-------------|-----------|----|----------|
|-----------|----------------|------------|-------------|-----------------|-------------|-------------|-----------|----|----------|

1 = This parameter is not certified by the primary accrediting authority (NY NELAC) for this matrix. NY NELAC does not offer certification for all parameters at this time.

RL/PQL=Reporting/Practical Quantitation Level (Equivalent to NELAC LOQ, Limit of Quantitation) ND=Not Detected BRL=Below Reporting Level L=Biased Low LOD=Limit of Detection MDL=Method Detection Limit1

QA/QC Surrogates: Surrogates are compounds (preceeded with a %) added by the lab to determine analysis efficiency. Surrogate results(%) listed in the report are not "detected" compounds.

Comments:

E = Estimated value quantitated above calibration range for this compound.

If you are the client above and have any questions concerning this testing, please do not hesitate to contact Phoenix Client Services at ext.200. The contents of this report cannot be discussed with anyone other than the client listed above without their written consent.

Phyllis Shiller, Laboratory Director

January 10, 2024

Reviewed and Released by: Greg Lawrence, Assistant Lab Director



Environmental Laboratories, Inc.
 587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045
 Tel. (860) 645-1102 Fax (860) 645-0823



Analysis Report

January 10, 2024

FOR: Attn:
 Advanced Cleanup Technologies, Inc.
 228 Park Ave S PMB 34864
 New York, New York 10003

Sample Information

Matrix: AIR
 Location Code: ACT
 Rush Request: Standard
 P.O.#:
 Canister Id: 23330
 Project ID: 9268-BXNY
 Client ID: IA-7

Custody Information

Collected by: YY,RB
 Received by: CP
 Analyzed by: see "By" below

Date Time
 08/02/22 10:22
 08/03/22 17:50

Laboratory Data

SDG ID: GCL97021
 Phoenix ID: CL97022

| Parameter | ppbv Result | ppbv RL | LOD/ MDL | ug/m3 Result | ug/m3 RL | LOD/ MDL | Date/Time | By | Dilution |
|-------------------------------|-------------|---------|----------|--------------|----------|----------|-----------|-----|----------|
| Volatiles (TO15) | | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | ND | 0.146 | 0.146 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,1,1-Trichloroethane | ND | 0.183 | 0.183 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,1,2,2-Tetrachloroethane | ND | 0.146 | 0.146 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,1,2-Trichloroethane | ND | 0.183 | 0.183 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,1-Dichloroethane | ND | 0.247 | 0.247 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,1-Dichloroethene | ND | 0.051 | 0.051 | ND | 0.20 | 0.20 | 08/03/22 | KCA | 1 |
| 1,2,4-Trichlorobenzene | ND | 0.135 | 0.135 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,2,4-Trimethylbenzene | 1.23 | 0.204 | 0.204 | 6.04 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,2-Dibromoethane(EDB) | ND | 0.130 | 0.130 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,2-Dichlorobenzene | ND | 0.166 | 0.166 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,2-Dichloroethane | ND | 0.247 | 0.247 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,2-dichloropropane | ND | 0.217 | 0.217 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,2-Dichlorotetrafluoroethane | ND | 0.143 | 0.143 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,3,5-Trimethylbenzene | 0.406 | 0.204 | 0.204 | 1.99 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,3-Butadiene | ND | 0.452 | 0.452 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,3-Dichlorobenzene | ND | 0.166 | 0.166 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,4-Dichlorobenzene | ND | 0.166 | 0.166 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 1,4-Dioxane | ND | 0.278 | 0.278 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 2-Hexanone(MBK) | ND | 0.244 | 0.244 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 4-Ethyltoluene | 1.03 | 0.204 | 0.204 | 5.06 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 4-Isopropyltoluene | ND | 0.182 | 0.182 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| 4-Methyl-2-pentanone(MIBK) | ND | 0.244 | 0.244 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Acetone | 44.7 | E 0.421 | 0.421 | 106 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Acrylonitrile | ND | 0.461 | 0.461 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Benzene | 1.72 | 0.313 | 0.313 | 5.49 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Benzyl chloride | ND | 0.193 | 0.193 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |

| Parameter | ppbv Result | ppbv RL | LOD/ MDL | ug/m3 Result | ug/m3 RL | LOD/ MDL | Date/Time | By | Dilution |
|--|----------------|------------|-------------|-----------------|-------------|-------------|-----------|-----|----------|
| Bromodichloromethane | ND | 0.149 | 0.149 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Bromoform | ND | 0.097 | 0.097 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Bromomethane | ND | 0.258 | 0.258 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Carbon Disulfide | ND | 0.321 | 0.321 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Carbon Tetrachloride | 0.081 | 0.032 | 0.032 | 0.51 | 0.20 | 0.20 | 08/03/22 | KCA | 1 |
| Chlorobenzene | ND | 0.217 | 0.217 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Chloroethane | ND | 0.379 | 0.379 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Chloroform | ND | 0.205 | 0.205 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Chloromethane | 0.630 | 0.485 | 0.485 | 1.30 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Cis-1,2-Dichloroethene | 16.0 | 0.051 | 0.051 | 63.4 | 0.20 | 0.20 | 08/03/22 | KCA | 1 |
| cis-1,3-Dichloropropene | ND | 0.221 | 0.221 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Cyclohexane | 2.16 | 0.291 | 0.291 | 7.43 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Dibromochloromethane | ND | 0.118 | 0.118 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Dichlorodifluoromethane | 0.502 | 0.202 | 0.202 | 2.48 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Ethanol | 76.1 | E 0.531 | 0.531 | 143 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Ethyl acetate | ND | 0.278 | 0.278 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Ethylbenzene | 0.995 | 0.230 | 0.230 | 4.32 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Heptane | 1.99 | 0.244 | 0.244 | 8.15 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Hexachlorobutadiene | ND | 0.094 | 0.094 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Hexane | 6.22 | 0.284 | 0.284 | 21.9 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Isopropylalcohol | 4.24 | 0.407 | 0.407 | 10.4 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Isopropylbenzene | ND | 0.204 | 0.204 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| m,p-Xylene | 3.54 | 0.230 | 0.230 | 15.4 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Methyl Ethyl Ketone | 1.17 | 0.339 | 0.339 | 3.45 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Methyl tert-butyl ether(MTBE) | ND | 0.278 | 0.278 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Methylene Chloride | ND | 0.863 | 0.863 | ND | 3.00 | 3.00 | 08/03/22 | KCA | 1 |
| n-Butylbenzene | ND | 0.182 | 0.182 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| o-Xylene | 1.33 | 0.230 | 0.230 | 5.77 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Propylene | 1.01 | 0.581 | 0.581 | 1.74 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| sec-Butylbenzene | ND | 0.182 | 0.182 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Styrene | ND | 0.235 | 0.235 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Tetrachloroethene | 2.80 | 0.037 | 0.037 | 19.0 | 0.25 | 0.25 | 08/03/22 | KCA | 1 |
| Tetrahydrofuran | 2.57 | 0.339 | 0.339 | 7.57 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Toluene | 7.10 | 0.266 | 0.266 | 26.7 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Trans-1,2-Dichloroethene | ND | 0.252 | 0.252 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| trans-1,3-Dichloropropene | ND | 0.221 | 0.221 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Trichloroethene | 1.90 | 0.037 | 0.037 | 10.2 | 0.20 | 0.20 | 08/03/22 | KCA | 1 |
| Trichlorofluoromethane | 0.289 | 0.178 | 0.178 | 1.62 | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Trichlorotrifluoroethane | ND | 0.131 | 0.131 | ND | 1.00 | 1.00 | 08/03/22 | KCA | 1 |
| Vinyl Chloride | 1.90 | 0.078 | 0.078 | 4.85 | 0.20 | 0.20 | 08/03/22 | KCA | 1 |
| <u>QA/QC Surrogates/Internals</u> | | | | | | | | | |
| % Bromofluorobenzene | 98 | % | % | 98 | % | % | 08/03/22 | KCA | 1 |
| % IS-1,4-Difluorobenzene | 108 | % | % | 108 | % | % | 08/03/22 | KCA | 1 |
| % IS-Bromochloromethane | 104 | % | % | 104 | % | % | 08/03/22 | KCA | 1 |
| % IS-Chlorobenzene-d5 | 109 | % | % | 109 | % | % | 08/03/22 | KCA | 1 |

| Parameter | ppbv Result | ppbv RL | LOD/ MDL | ug/m3 Result | ug/m3LOD/ RL MDL | Date/Time | By | Dilution |
|-----------|----------------|------------|-------------|-----------------|---------------------|-----------|----|----------|
|-----------|----------------|------------|-------------|-----------------|---------------------|-----------|----|----------|

1 = This parameter is not certified by the primary accrediting authority (NY NELAC) for this matrix. NY NELAC does not offer certification for all parameters at this time.

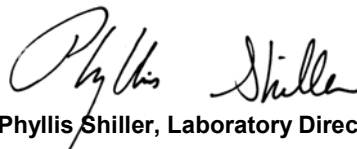
RL/PQL=Reporting/Practical Quantitation Level (Equivalent to NELAC LOQ, Limit of Quantitation) ND=Not Detected BRL=Below Reporting Level L=Biased Low LOD=Limit of Detection MDL=Method Detection Limit1

QA/QC Surrogates: Surrogates are compounds (preceeded with a %) added by the lab to determine analysis efficiency. Surrogate results(%) listed in the report are not "detected" compounds.

Comments:

E = Estimated value quantitated above calibration range for this compound.

If you are the client above and have any questions concerning this testing, please do not hesitate to contact Phoenix Client Services at ext.200. The contents of this report cannot be discussed with anyone other than the client listed above without their written consent.



Phyllis Shiller, Laboratory Director

January 10, 2024

Reviewed and Released by: Greg Lawrence, Assistant Lab Director



Environmental Laboratories, Inc.
 587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045
 Tel. (860) 645-1102 Fax (860) 645-0823



Canister Sampling Information

January 10, 2024

FOR: Attn:
 Advanced Cleanup Technologies, Inc.
 228 Park Ave S PMB 34864
 New York, New York 10003

Location Code: ACT

SDG I.D.: GCL97021

Project ID: 9268-BXNY

| Client Id | Lab Id | Canister | | Reg. Id | Chk Out Date | Laboratory | | | | | Field | | | |
|-----------|---------|----------|------|---------|--------------|------------|-------|----------|---------|----------|----------|--------|---------------------|-------------------|
| | | Id | Type | | | Out Hg | In Hg | Out Flow | In Flow | Flow RPD | Start Hg | End Hg | Sampling Start Date | Sampling End Date |
| IA-4 | CL97021 | 28560 | 6.0L | 5043 | 07/19/22 | -30 | -7 | 3.6 | 4 | 10.5 | -30 | -8 | 08/01/22 15:00 | 08/02/22 11:03 |
| IA-7 | CL97022 | 23330 | 6.0L | 5000 | 07/19/22 | -30 | -5 | 3.8 | 4.2 | 10.0 | -30 | -6 | 08/01/22 15:00 | 08/02/22 10:22 |



Environmental Laboratories, Inc.
587 East Middle Turnpike, P.O.Box 370, Manchester, CT 06045
Tel. (860) 645-1102



QA/QC Report

January 10, 2024

QA/QC Data

SDG I.D.: GCL97021

| Parameter | Blk ppbv | Blk RL ppbv | Blk ug/m3 | Blk RL ug/m3 | LCS % | Sample Result ug/m3 | Sample Dup ug/m3 | Sample Result ppbv | Sample Dup ppbv | DUP RPD | % Rec Limits | % RPD Limits |
|---|-------------|-------------------|--------------|--------------------|----------|---------------------------|------------------------|--------------------------|-----------------------|------------|--------------------|--------------------|
| QA/QC Batch 635992 (ppbv), QC Sample No: CL97022 (CL97021, CL97022) | | | | | | | | | | | | |
| Volatiles | | | | | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | ND | 0.150 | ND | 1.03 | 113 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,1,1-Trichloroethane | ND | 0.180 | ND | 0.98 | 104 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,1,2,2-Tetrachloroethane | ND | 0.150 | ND | 1.03 | 97 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,1,2-Trichloroethane | ND | 0.180 | ND | 0.98 | 100 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,1-Dichloroethane | ND | 0.250 | ND | 1.01 | 106 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,1-Dichloroethene | ND | 0.050 | ND | 0.20 | 110 | ND | 0.22 | ND | 0.056 | NC | 70 - 130 | 25 |
| 1,2,4-Trichlorobenzene | ND | 0.130 | ND | 0.96 | 149 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,2,4-Trimethylbenzene | ND | 0.200 | ND | 0.98 | 98 | 6.04 | 6.19 | 1.23 | 1.26 | 2.4 | 70 - 130 | 25 |
| 1,2-Dibromoethane(EDB) | ND | 0.130 | ND | 1.00 | 102 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,2-Dichlorobenzene | ND | 0.170 | ND | 1.02 | 101 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,2-Dichloroethane | ND | 0.250 | ND | 1.01 | 110 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,2-dichloropropane | ND | 0.220 | ND | 1.02 | 104 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,2-Dichlorotetrafluoroethane | ND | 0.140 | ND | 0.98 | 107 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,3,5-Trimethylbenzene | ND | 0.200 | ND | 0.98 | 95 | 1.99 | 1.98 | 0.406 | 0.404 | NC | 70 - 130 | 25 |
| 1,3-Butadiene | ND | 0.450 | ND | 0.99 | 120 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,3-Dichlorobenzene | ND | 0.170 | ND | 1.02 | 105 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,4-Dichlorobenzene | ND | 0.170 | ND | 1.02 | 103 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 1,4-Dioxane | ND | 0.280 | ND | 1.01 | 96 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 2-Hexanone(MBK) | ND | 0.240 | ND | 0.98 | 128 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 4-Ethyltoluene | ND | 0.200 | ND | 0.98 | 100 | 5.06 | 5.01 | 1.03 | 1.02 | 1.0 | 70 - 130 | 25 |
| 4-Isopropyltoluene | ND | 0.180 | ND | 0.99 | 99 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| 4-Methyl-2-pentanone(MIBK) | ND | 0.240 | ND | 0.98 | 128 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Acetone | ND | 0.420 | ND | 1.00 | 132 | 106 E | 107 | 44.7 E | 44.9 | 0.4 | 70 - 130 | 25 |
| Acrylonitrile | ND | 0.460 | ND | 1.00 | 129 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Benzene | ND | 0.310 | ND | 0.99 | 98 | 5.49 | 5.49 | 1.72 | 1.72 | 0.0 | 70 - 130 | 25 |
| Benzyl chloride | ND | 0.190 | ND | 0.98 | 110 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Bromodichloromethane | ND | 0.150 | ND | 1.00 | 111 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Bromoform | ND | 0.097 | ND | 1.00 | 112 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Bromomethane | ND | 0.260 | ND | 1.01 | 98 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Carbon Disulfide | ND | 0.320 | ND | 1.00 | 105 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Carbon Tetrachloride | ND | 0.032 | ND | 0.20 | 112 | 0.51 | 0.52 | 0.081 | 0.083 | NC | 70 - 130 | 25 |
| Chlorobenzene | ND | 0.220 | ND | 1.01 | 96 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Chloroethane | ND | 0.380 | ND | 1.00 | 93 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Chloroform | ND | 0.200 | ND | 0.98 | 102 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Chloromethane | ND | 0.480 | ND | 0.99 | 134 | 1.30 | 1.40 | 0.630 | 0.676 | NC | 70 - 130 | 25 |
| Cis-1,2-Dichloroethene | ND | 0.050 | ND | 0.20 | 117 | 63.4 | 62.6 | 16.0 | 15.8 | 1.3 | 70 - 130 | 25 |
| cis-1,3-Dichloropropene | ND | 0.220 | ND | 1.00 | 102 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Cyclohexane | ND | 0.290 | ND | 1.00 | 99 | 7.43 | 7.67 | 2.16 | 2.23 | 3.2 | 70 - 130 | 25 |
| Dibromochloromethane | ND | 0.120 | ND | 1.02 | 110 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Dichlorodifluoromethane | ND | 0.200 | ND | 0.99 | 107 | 2.48 | 2.57 | 0.502 | 0.520 | NC | 70 - 130 | 25 |
| Ethanol | ND | 0.530 | ND | 1.00 | 149 | 143 E | 106 | 76.1 E | 56.4 | 29.7 | 70 - 130 | 25 |

QA/QC Data

SDG I.D.: GCL97021


| Parameter | Bik ppbv | Bik RL ppbv | Bik ug/m3 | Bik RL ug/m3 | LCS % | Sample Result ug/m3 | Sample Dup ug/m3 | Sample Result ppbv | Sample Dup ppbv | DUP RPD | % Rec Limits | % RPD Limits |
|-------------------------------|-------------|-------------------|--------------|--------------------|----------|---------------------------|------------------------|--------------------------|-----------------------|------------|--------------------|--------------------|
| Ethyl acetate | ND | 0.280 | ND | 1.01 | 94 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Ethylbenzene | ND | 0.230 | ND | 1.00 | 92 | 4.32 | 4.29 | 0.995 | 0.989 | NC | 70 - 130 | 25 |
| Heptane | ND | 0.240 | ND | 0.98 | 126 | 8.15 | 8.15 | 1.99 | 1.99 | 0.0 | 70 - 130 | 25 |
| Hexachlorobutadiene | ND | 0.094 | ND | 1.00 | 126 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Hexane | ND | 0.280 | ND | 0.99 | 111 | 21.9 | 22.2 | 6.22 | 6.30 | 1.3 | 70 - 130 | 25 |
| Isopropylalcohol | ND | 0.410 | ND | 1.01 | 130 | 10.4 | 8.72 | 4.24 | 3.55 | 17.7 | 70 - 130 | 25 |
| Isopropylbenzene | ND | 0.200 | ND | 0.98 | 106 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| m,p-Xylene | ND | 0.230 | ND | 1.00 | 95 | 15.4 | 15.7 | 3.54 | 3.62 | 2.2 | 70 - 130 | 25 |
| Methyl Ethyl Ketone | ND | 0.340 | ND | 1.00 | 129 | 3.45 | 3.65 | 1.17 | 1.24 | NC | 70 - 130 | 25 |
| Methyl tert-butyl ether(MTBE) | ND | 0.280 | ND | 1.01 | 97 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Methylene Chloride | ND | 0.860 | ND | 2.99 | 125 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| n-Butylbenzene | ND | 0.180 | ND | 0.99 | 102 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| o-Xylene | ND | 0.230 | ND | 1.00 | 94 | 5.77 | 5.95 | 1.33 | 1.37 | 3.0 | 70 - 130 | 25 |
| Propylene | ND | 0.580 | ND | 1.00 | 98 | 1.74 | 1.87 | 1.01 | 1.09 | NC | 70 - 130 | 25 |
| sec-Butylbenzene | ND | 0.180 | ND | 0.99 | 101 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Styrene | ND | 0.230 | ND | 0.98 | 96 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Tetrachloroethene | ND | 0.037 | ND | 0.25 | 99 | 19.0 | 19.0 | 2.80 | 2.80 | 0.0 | 70 - 130 | 25 |
| Tetrahydrofuran | ND | 0.340 | ND | 1.00 | 128 | 7.57 | 7.75 | 2.57 | 2.63 | 2.3 | 70 - 130 | 25 |
| Toluene | ND | 0.270 | ND | 1.02 | 100 | 26.7 | 26.9 | 7.10 | 7.15 | 0.7 | 70 - 130 | 25 |
| Trans-1,2-Dichloroethene | ND | 0.250 | ND | 0.99 | 108 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| trans-1,3-Dichloropropene | ND | 0.220 | ND | 1.00 | 102 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Trichloroethene | ND | 0.037 | ND | 0.20 | 102 | 10.2 | 10.9 | 1.90 | 2.03 | 6.6 | 70 - 130 | 25 |
| Trichlorofluoromethane | ND | 0.180 | ND | 1.01 | 107 | 1.62 | 1.59 | 0.289 | 0.284 | NC | 70 - 130 | 25 |
| Trichlorotrifluoroethane | ND | 0.130 | ND | 1.00 | 107 | ND | ND | ND | ND | NC | 70 - 130 | 25 |
| Vinyl Chloride | ND | 0.078 | ND | 0.20 | 108 | 4.85 | 4.80 | 1.90 | 1.88 | 1.1 | 70 - 130 | 25 |
| % Bromofluorobenzene | 96 | % | 96 | % | 99 | 98 | 99 | 98 | 99 | NC | 70 - 130 | 25 |
| % IS-1,4-Difluorobenzene | 111 | % | 111 | % | 110 | 108 | 105 | 108 | 105 | NC | 60 - 140 | 25 |
| % IS-Bromochloromethane | 109 | % | 109 | % | 106 | 104 | 103 | 104 | 103 | NC | 60 - 140 | 25 |
| % IS-Chlorobenzene-d5 | 108 | % | 108 | % | 119 | 109 | 106 | 109 | 106 | NC | 60 - 140 | 25 |

l = This parameter is outside laboratory LCS/LCSD specified recovery limits.

r = This parameter is outside laboratory RPD specified recovery limits.

If there are any questions regarding this data, please call Phoenix Client Services at extension 200.

- RPD - Relative Percent Difference
- LCS - Laboratory Control Sample
- LCSD - Laboratory Control Sample Duplicate
- MS - Matrix Spike
- MS Dup - Matrix Spike Duplicate
- NC - No Criteria
- Intf - Interference


 Phyllis Shiller, Laboratory Director
 January 10, 2024

Wednesday, January 10, 2024

Criteria: None

State: NY

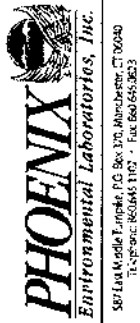
Sample Criteria Exceedances Report

GCL97021 - ACT

| SampNo | Acode | Phoenix Analyte | Criteria | Result | RL | Criteria | RL Criteria | Analysis Units |
|--------|-------|-----------------|----------|--------|----|----------|----------------|-------------------|
|--------|-------|-----------------|----------|--------|----|----------|----------------|-------------------|

*** No Data to Display ***

Phoenix Laboratories does not assume responsibility for the data contained in this exceedance report. It is provided as an additional tool to identify requested criteria exceedences. All efforts are made to ensure the accuracy of the data (obtained from appropriate agencies). A lack of exceedence information does not necessarily suggest conformance to the criteria. It is ultimately the site professional's responsibility to determine appropriate compliance.



CHAIN OF CUSTODY RECORD
AIR ANALYSES

5811 Middlebrook, D.C. Box 370, Bethesda, CT 06040
Telephone: 860/665-1102 • Fax: 860/665-3623

860-645-1102
email: greg@phoenixlabs.com

ACT

Report to: **Jason Stewart**
 Customer: **Advanced Cleanup Technologies, Inc.**
 Address: **228 Park Ave S PMB 34864**
13763 New York, New York 10003

Project Name: **q268 - Bx NY**
 Invoice to: **Karen Friedman**
 Requested Deliverable: **RCP ASP CAT B**
 Other: **MCP NJ Deliverables**

Requested Deliverable: **RCP ASP CAT B**
 Quote Number: _____

| Phoenix ID # | Client Sample ID | Canister ID # | Canister Size (L) | Outgoing Canister Pressure (Psi) | Incoming Canister Pressure (Psi) | Flow Regulator ID # | Flow Controller Setting (mL/min) | THIS SECTION FOR LAB USE ONLY | | | | Canister Pressure at End (Psi) | Canister Pressure at Start (Psi) | Sample Start Date | Sampling End Time | Sampling Start Time | Ambient/Indoor Air | Soil Gas | Grab (G) Composite (C) | TO-15 | ANALYSES |
|--------------|------------------|---------------|-------------------|----------------------------------|----------------------------------|---------------------|----------------------------------|-------------------------------|-------------------|----------------------------------|----------------------------------|--------------------------------|----------------------------------|-------------------|-------------------|---------------------|--------------------|----------|------------------------|-------|----------|
| | | | | | | | | Canister ID # | Canister Size (L) | Outgoing Canister Pressure (Psi) | Incoming Canister Pressure (Psi) | | | | | | | | | | |
| 97021 | IA-4 | 28560 | 6.0L | -30 | MVA | 5043 | 3.6 | 15:00 8/1/22 | 11:05 8/2/22 | 8/1/22 | -30 | -8 | 8/1/22 | 8/1/22 | 8/1/22 | 8/1/22 | V | | | V | |
| 97022 | IA-7 | 23338 | 6.0L | -30 | MVA | 5042 | 3.6 | 15:00 8/1/22 | 10:27 8/1/22 | 8/1/22 | -30 | -6 | 8/1/22 | 8/1/22 | 8/1/22 | 8/1/22 | V | | | V | |
| | | 23330 | 6.0L | -30 | MVA | 5000 | 3.8 | | | | | | | | | | | | | | |

Relinquished by: Ray Buzeta 8/2/22 3:15 PM
 Accepted by: Greg Friedman 8-3-22 10:54
 Date: 8-3-22
 Time: 10:54

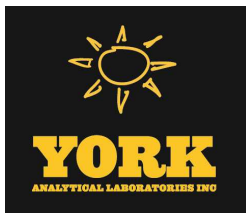
Signature: _____ Date: _____

Requested Criteria: (Please Circle)

| | | | | |
|-------------|------------------------|----------------------------|----------------------------|----------------------------|
| CT: TAC I/C | Indoor Air Residential | NI: Indoor Air Residential | PA: Indoor Air Residential | VT: Indoor Air Residential |
| TAC RES | Ind/Commercial | Ind/Commercial | Residential | Residential |
| SVVC I/C | Soil Gas | Soil Gas | Non-residential | Industrial |
| SVVC RES | Residential | Residential | | |
| GWV I/C | Ind/Commercial | Ind/Commercial | | |
| GWV RES | | | | |

Turnaround Time: 1 Day 2 Day 3 Day 4 Day 5 Day

SPECIAL INSTRUCTIONS, OC REQUIREMENTS, REGULATORY INFORMATION:
 Can # 23338 (3) - 6.0L 24 hr returned unused



Technical Report

prepared for:

Advanced Cleanup Technologies, Inc.
228 Park Ave S PMB 34864
New York NY, 10003
Attention: Paul Stewart

Report Date: 03/22/2023
Client Project ID: 9628-BXNY
York Project (SDG) No.: 23C0637

CT Cert. No. PH-0723

New Jersey Cert. No. CT005 and NY037



New York Cert. Nos. 10854 and 12058

PA Cert. No. 68-04440

120 RESEARCH DRIVE
www.YORKLAB.com

STRATFORD, CT 06615
(203) 325-1371



132-02 89th AVENUE
FAX (203) 357-0166

RICHMOND HILL, NY 11418
ClientServices@yorklab.com

Report Date: 03/22/2023
Client Project ID: 9628-BXNY
York Project (SDG) No.: 23C0637

Advanced Cleanup Technologies, Inc.
228 Park Ave S PMB 34864
New York NY, 10003
Attention: Paul Stewart

Purpose and Results

This report contains the analytical data for the sample(s) identified on the attached chain-of-custody received in our laboratory on March 10, 2023 and listed below. The project was identified as your project: **9628-BXNY**.

The analyses were conducted utilizing appropriate EPA, Standard Methods, and ASTM methods as detailed in the data summary tables.

All samples were received in proper condition meeting the customary acceptance requirements for environmental samples except those indicated under the Sample and Analysis Qualifiers section of this report.

All analyses met the method and laboratory standard operating procedure requirements except as indicated by any data flags, the meaning of which are explained in the Sample and Data Qualifiers Relating to This Work Order section of this report and case narrative if applicable.

The results of the analyses, which are all reported on dry weight basis (soils) unless otherwise noted, are detailed in the following pages.

Please contact Client Services at 203.325.1371 with any questions regarding this report.

| <u>York Sample ID</u> | <u>Client Sample ID</u> | <u>Matrix</u> | <u>Date Collected</u> | <u>Date Received</u> |
|-----------------------|-------------------------|---------------|-----------------------|----------------------|
| 23C0637-01 | DW-1 | Water | 03/09/2023 | 03/10/2023 |
| 23C0637-02 | DW-2 | Water | 03/09/2023 | 03/10/2023 |

General Notes for York Project (SDG) No.: 23C0637

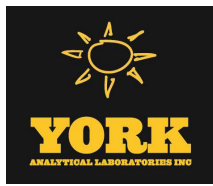
1. The RLs and MDLs (Reporting Limit and Method Detection Limit respectively) reported are adjusted for any dilution necessary due to the levels of target and/or non-target analytes and matrix interference. The RL(REPORTING LIMIT) is based upon the lowest standard utilized for the calibration where applicable.
2. Samples are retained for a period of thirty days after submittal of report, unless other arrangements are made.
3. York's liability for the above data is limited to the dollar value paid to York for the referenced project.
4. This report shall not be reproduced without the written approval of York Analytical Laboratories, Inc.
5. All analyses conducted met method or Laboratory SOP requirements. See the Sample and Data Qualifiers Section for further information.
6. It is noted that no analyses reported herein were subcontracted to another laboratory, unless noted in the report.
7. This report reflects results that relate only to the samples submitted on the attached chain-of-custody form(s) received by York.
8. Analyses conducted at York Analytical Laboratories, Inc. Stratford, CT are indicated by NY Cert. No. 10854; those conducted at York Analytical Laboratories, Inc., Richmond Hill, NY are indicated by NY Cert. No. 12058.

Approved By: 

Date: 03/22/2023

Cassie L. Mosher
Laboratory Manager





Sample Information

Client Sample ID: DW-1

York Sample ID: 23C0637-01

York Project (SDG) No.

Client Project ID

Matrix

Collection Date/Time

Date Received

23C0637

9628-BXNY

Water

March 9, 2023 1:20 pm

03/10/2023

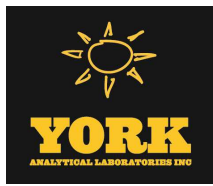
VOA, 8260 LOW MASTER

Log-in Notes:

Sample Notes:

Sample Prepared by Method: EPA 5030B

| CAS No. | Parameter | Result | Flag | Units | Reported to LOD/MDL | LOQ | Dilution | Reference Method | Date/Time Prepared | Date/Time Analyzed | Analyst |
|----------|---|-------------|------|-------|---------------------|-------|----------|--|--------------------|--------------------|---------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | ND | | ug/L | 0.216 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 71-55-6 | 1,1,1-Trichloroethane | ND | | ug/L | 0.266 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | | ug/L | 0.256 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | ND | | ug/L | 0.286 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 79-00-5 | 1,1,2-Trichloroethane | ND | | ug/L | 0.249 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 75-34-3 | 1,1-Dichloroethane | ND | | ug/L | 0.272 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 75-35-4 | 1,1-Dichloroethylene | ND | | ug/L | 0.327 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | | ug/L | 0.222 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 96-18-4 | 1,2,3-Trichloropropane | ND | | ug/L | 0.273 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | | ug/L | 0.138 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 95-63-6 | 1,2,4-Trimethylbenzene | ND | | ug/L | 0.310 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | | ug/L | 0.432 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 106-93-4 | 1,2-Dibromoethane | ND | | ug/L | 0.215 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 95-50-1 | 1,2-Dichlorobenzene | ND | | ug/L | 0.270 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 107-06-2 | 1,2-Dichloroethane | ND | | ug/L | 0.377 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 78-87-5 | 1,2-Dichloropropane | ND | | ug/L | 0.327 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 108-67-8 | 1,3,5-Trimethylbenzene | ND | | ug/L | 0.347 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 541-73-1 | 1,3-Dichlorobenzene | ND | | ug/L | 0.283 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 106-46-7 | 1,4-Dichlorobenzene | ND | | ug/L | 0.311 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 123-91-1 | 1,4-Dioxane | ND | | ug/L | 35.3 | 80.0 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 78-93-3 | 2-Butanone | 11.3 | | ug/L | 0.421 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 591-78-6 | 2-Hexanone | ND | | ug/L | 0.320 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |



Sample Information

Client Sample ID: DW-1

York Sample ID: 23C0637-01

York Project (SDG) No.

Client Project ID

Matrix

Collection Date/Time

Date Received

23C0637

9628-BXNY

Water

March 9, 2023 1:20 pm

03/10/2023

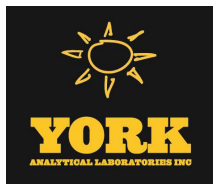
VOA, 8260 LOW MASTER

Log-in Notes:

Sample Notes:

Sample Prepared by Method: EPA 5030B

| CAS No. | Parameter | Result | Flag | Units | Reported to LOD/MDL | LOQ | Dilution | Reference Method | Date/Time Prepared | Date/Time Analyzed | Analyst |
|------------|---------------------------|--------|-------|-------|---------------------|-------|----------|--|--------------------|--------------------|---------|
| 108-10-1 | 4-Methyl-2-pentanone | ND | | ug/L | 0.365 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 67-64-1 | Acetone | 9.72 | | ug/L | 1.34 | 2.00 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 107-02-8 | Acrolein | ND | | ug/L | 0.447 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 107-13-1 | Acrylonitrile | ND | | ug/L | 0.422 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 71-43-2 | Benzene | ND | | ug/L | 0.279 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 74-97-5 | Bromochloromethane | ND | | ug/L | 0.354 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 75-27-4 | Bromodichloromethane | ND | | ug/L | 0.245 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 75-25-2 | Bromoform | ND | CCVE | ug/L | 0.163 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 74-83-9 | Bromomethane | ND | | ug/L | 0.119 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 75-15-0 | Carbon disulfide | ND | | ug/L | 0.362 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 56-23-5 | Carbon tetrachloride | ND | | ug/L | 0.204 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 108-90-7 | Chlorobenzene | ND | | ug/L | 0.284 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 75-00-3 | Chloroethane | ND | | ug/L | 0.448 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 67-66-3 | Chloroform | 2.90 | | ug/L | 0.243 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 74-87-3 | Chloromethane | ND | | ug/L | 0.372 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 156-59-2 | cis-1,2-Dichloroethylene | 1.37 | | ug/L | 0.294 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 10061-01-5 | cis-1,3-Dichloropropylene | ND | | ug/L | 0.262 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 110-82-7 | Cyclohexane | ND | QL-02 | ug/L | 0.491 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 124-48-1 | Dibromochloromethane | ND | | ug/L | 0.146 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 74-95-3 | Dibromomethane | ND | | ug/L | 0.203 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 75-71-8 | Dichlorodifluoromethane | ND | | ug/L | 0.451 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 100-41-4 | Ethyl Benzene | ND | | ug/L | 0.290 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 87-68-3 | Hexachlorobutadiene | ND | CCVE | ug/L | 0.241 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |



Sample Information

Client Sample ID: DW-1

York Sample ID: 23C0637-01

York Project (SDG) No.

Client Project ID

Matrix

Collection Date/Time

Date Received

23C0637

9628-BXNY

Water

March 9, 2023 1:20 pm

03/10/2023

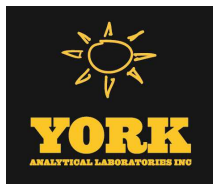
VOA, 8260 LOW MASTER

Log-in Notes:

Sample Notes:

Sample Prepared by Method: EPA 5030B

| CAS No. | Parameter | Result | Flag | Units | Reported to LOD/MDL | LOQ | Dilution | Reference Method | Date/Time Prepared | Date/Time Analyzed | Analyst |
|-------------|---------------------------------------|---------------|---------------|-------|-------------------------|-------|----------|--|--------------------|--------------------|---------|
| 98-82-8 | Isopropylbenzene | ND | | ug/L | 0.405 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 79-20-9 | Methyl acetate | ND | | ug/L | 0.442 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 0.330 | J | ug/L | 0.244 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PA | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 108-87-2 | Methylcyclohexane | ND | | ug/L | 0.477 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 75-09-2 | Methylene chloride | ND | | ug/L | 0.397 | 2.00 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 104-51-8 | n-Butylbenzene | ND | | ug/L | 0.399 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 103-65-1 | n-Propylbenzene | ND | | ug/L | 0.384 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 95-47-6 | o-Xylene | ND | | ug/L | 0.261 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 179601-23-1 | p- & m- Xylenes | ND | | ug/L | 0.578 | 1.00 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 99-87-6 | p-Isopropyltoluene | ND | | ug/L | 0.377 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 135-98-8 | sec-Butylbenzene | ND | | ug/L | 0.444 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 100-42-5 | Styrene | ND | | ug/L | 0.255 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 75-65-0 | tert-Butyl alcohol (TBA) | 1.17 | CCVE, ICVE | ug/L | 0.608 | 1.00 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 98-06-6 | tert-Butylbenzene | ND | | ug/L | 0.367 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 127-18-4 | Tetrachloroethylene | 5.95 | | ug/L | 0.239 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PA | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 108-88-3 | Toluene | 1.64 | | ug/L | 0.346 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PA | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 156-60-5 | trans-1,2-Dichloroethylene | ND | | ug/L | 0.279 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 10061-02-6 | trans-1,3-Dichloropropylene | ND | | ug/L | 0.229 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 79-01-6 | Trichloroethylene | ND | | ug/L | 0.249 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 75-69-4 | Trichlorofluoromethane | ND | | ug/L | 0.337 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 75-01-4 | Vinyl Chloride | ND | | ug/L | 0.469 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| 1330-20-7 | Xylenes, Total | ND | | ug/L | 0.836 | 1.50 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP | 03/20/2023 06:45 | 03/21/2023 09:04 | JTG |
| | Surrogate Recoveries | Result | | | Acceptance Range | | | | | | |



Sample Information

Client Sample ID: DW-1

York Sample ID: 23C0637-01

York Project (SDG) No.

Client Project ID

Matrix

Collection Date/Time

Date Received

23C0637

9628-BXNY

Water

March 9, 2023 1:20 pm

03/10/2023

VOA, 8260 LOW MASTER

Log-in Notes:

Sample Notes:

Sample Prepared by Method: EPA 5030B

| CAS No. | Parameter | Result | Flag | Units | Reported to LOD/MDL | LOQ | Dilution | Reference Method | Date/Time Prepared | Date/Time Analyzed | Analyst |
|------------|--|--------|------|-------|---------------------|-----|----------|------------------|--------------------|--------------------|---------|
| 17060-07-0 | Surrogate: SURRE: 1,2-Dichloroethane-d4 | 103 % | | | 69-130 | | | | | | |
| 2037-26-5 | Surrogate: SURRE: Toluene-d8 | 101 % | | | 81-117 | | | | | | |
| 460-00-4 | Surrogate: SURRE: p-Bromofluorobenzene | 86.3 % | | | 79-122 | | | | | | |

Sample Information

Client Sample ID: DW-2

York Sample ID: 23C0637-02

York Project (SDG) No.

Client Project ID

Matrix

Collection Date/Time

Date Received

23C0637

9628-BXNY

Water

March 9, 2023 12:45 pm

03/10/2023

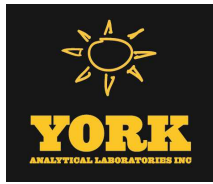
VOA, 8260 LOW MASTER

Log-in Notes:

Sample Notes:

Sample Prepared by Method: EPA 5030B

| CAS No. | Parameter | Result | Flag | Units | Reported to LOD/MDL | LOQ | Dilution | Reference Method | Date/Time Prepared | Date/Time Analyzed | Analyst |
|----------|---|-------------|------|-------|---------------------|-------|----------|--|--------------------|--------------------|---------|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | ND | | ug/L | 0.216 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 71-55-6 | 1,1,1-Trichloroethane | ND | | ug/L | 0.266 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | | ug/L | 0.256 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | ND | | ug/L | 0.286 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 79-00-5 | 1,1,2-Trichloroethane | ND | | ug/L | 0.249 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 75-34-3 | 1,1-Dichloroethane | ND | | ug/L | 0.272 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 75-35-4 | 1,1-Dichloroethylene | 61.8 | CCVE | ug/L | 0.327 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | | ug/L | 0.222 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 96-18-4 | 1,2,3-Trichloropropane | ND | | ug/L | 0.273 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | | ug/L | 0.138 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 95-63-6 | 1,2,4-Trimethylbenzene | ND | | ug/L | 0.310 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | | ug/L | 0.432 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 106-93-4 | 1,2-Dibromoethane | ND | | ug/L | 0.215 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 95-50-1 | 1,2-Dichlorobenzene | ND | | ug/L | 0.270 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |



Sample Information

Client Sample ID: DW-2

York Sample ID: 23C0637-02

York Project (SDG) No.

Client Project ID

Matrix

Collection Date/Time

Date Received

23C0637

9628-BXNY

Water

March 9, 2023 12:45 pm

03/10/2023

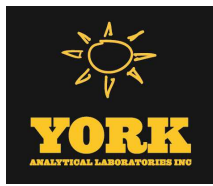
VOA, 8260 LOW MASTER

Log-in Notes:

Sample Notes:

Sample Prepared by Method: EPA 5030B

| CAS No. | Parameter | Result | Flag | Units | Reported to LOD/MDL | LOQ | Dilution | Reference Method | Date/Time Prepared | Date/Time Analyzed | Analyst |
|----------|------------------------|-------------|------|-------|---------------------|-------|----------|--|--------------------|--------------------|---------|
| 107-06-2 | 1,2-Dichloroethane | ND | | ug/L | 0.377 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 78-87-5 | 1,2-Dichloropropane | ND | | ug/L | 0.327 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 108-67-8 | 1,3,5-Trimethylbenzene | ND | | ug/L | 0.347 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 541-73-1 | 1,3-Dichlorobenzene | ND | | ug/L | 0.283 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 106-46-7 | 1,4-Dichlorobenzene | ND | | ug/L | 0.311 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 123-91-1 | 1,4-Dioxane | ND | | ug/L | 35.3 | 80.0 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 78-93-3 | 2-Butanone | ND | | ug/L | 0.421 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 591-78-6 | 2-Hexanone | ND | | ug/L | 0.320 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 108-10-1 | 4-Methyl-2-pentanone | ND | | ug/L | 0.365 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 67-64-1 | Acetone | 37.6 | | ug/L | 1.34 | 2.00 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 107-02-8 | Acrolein | ND | | ug/L | 0.447 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 107-13-1 | Acrylonitrile | ND | | ug/L | 0.422 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 71-43-2 | Benzene | 1.32 | | ug/L | 0.279 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 74-97-5 | Bromochloromethane | ND | | ug/L | 0.354 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 75-27-4 | Bromodichloromethane | ND | | ug/L | 0.245 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 75-25-2 | Bromoform | ND | CCVE | ug/L | 0.163 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 74-83-9 | Bromomethane | ND | | ug/L | 0.119 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 75-15-0 | Carbon disulfide | ND | | ug/L | 0.362 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 56-23-5 | Carbon tetrachloride | ND | | ug/L | 0.204 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 108-90-7 | Chlorobenzene | ND | | ug/L | 0.284 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 75-00-3 | Chloroethane | 9.22 | | ug/L | 0.448 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 67-66-3 | Chloroform | ND | | ug/L | 0.243 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 74-87-3 | Chloromethane | ND | | ug/L | 0.372 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |



Sample Information

Client Sample ID: DW-2

York Sample ID: 23C0637-02

York Project (SDG) No.

Client Project ID

Matrix

Collection Date/Time

Date Received

23C0637

9628-BXNY

Water

March 9, 2023 12:45 pm

03/10/2023

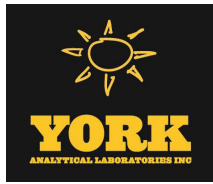
VOA, 8260 LOW MASTER

Log-in Notes:

Sample Notes:

Sample Prepared by Method: EPA 5030B

| CAS No. | Parameter | Result | Flag | Units | Reported to LOD/MDL | LOQ | Dilution | Reference Method | Date/Time Prepared | Date/Time Analyzed | Analyst |
|-------------|--------------------------------|--------|-------------------|-------|---------------------|-------|----------|--|--------------------|--------------------|---------|
| 156-59-2 | cis-1,2-Dichloroethylene | 22100 | | ug/L | 58.8 | 100 | 200 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/21/2023 06:13 | 03/22/2023 02:23 | JTG |
| 10061-01-5 | cis-1,3-Dichloropropylene | ND | | ug/L | 0.262 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 110-82-7 | Cyclohexane | 2.24 | CCVE, ICVE, QL-02 | ug/L | 0.491 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 124-48-1 | Dibromochloromethane | ND | | ug/L | 0.146 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 74-95-3 | Dibromomethane | ND | | ug/L | 0.203 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 75-71-8 | Dichlorodifluoromethane | ND | | ug/L | 0.451 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 100-41-4 | Ethyl Benzene | ND | | ug/L | 0.290 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 87-68-3 | Hexachlorobutadiene | ND | CCVE | ug/L | 0.241 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 98-82-8 | Isopropylbenzene | 3.41 | | ug/L | 0.405 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 79-20-9 | Methyl acetate | ND | | ug/L | 0.442 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | ND | | ug/L | 0.244 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 108-87-2 | Methylcyclohexane | 2.77 | CCVE | ug/L | 0.477 | 0.500 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 75-09-2 | Methylene chloride | ND | | ug/L | 0.397 | 2.00 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 104-51-8 | n-Butylbenzene | ND | | ug/L | 0.399 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 103-65-1 | n-Propylbenzene | ND | | ug/L | 0.384 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 95-47-6 | o-Xylene | 2.30 | | ug/L | 0.261 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 179601-23-1 | p- & m- Xylenes | ND | | ug/L | 0.578 | 1.00 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 99-87-6 | p-Isopropyltoluene | ND | | ug/L | 0.377 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 135-98-8 | sec-Butylbenzene | 0.870 | | ug/L | 0.444 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 100-42-5 | Styrene | ND | | ug/L | 0.255 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 75-65-0 | tert-Butyl alcohol (TBA) | ND | CCVE | ug/L | 0.608 | 1.00 | 1 | EPA 8260C Certifications: NELAC-NY10854,NELAC-NY12058,NJDEP,PADEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 98-06-6 | tert-Butylbenzene | ND | | ug/L | 0.367 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 127-18-4 | Tetrachloroethylene | 198 | | ug/L | 47.8 | 100 | 200 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PAI | 03/21/2023 06:13 | 03/22/2023 02:23 | JTG |



Sample Information

Client Sample ID: DW-2

York Sample ID: 23C0637-02

York Project (SDG) No.

Client Project ID

Matrix

Collection Date/Time

Date Received

23C0637

9628-BXNY

Water

March 9, 2023 12:45 pm

03/10/2023

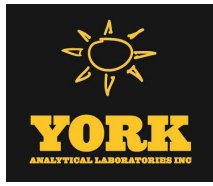
VOA, 8260 LOW MASTER

Log-in Notes:

Sample Notes:

Sample Prepared by Method: EPA 5030B

| CAS No. | Parameter | Result | Flag | Units | Reported to LOD/MDL | LOQ | Dilution | Reference Method | Date/Time Prepared | Date/Time Analyzed | Analyst |
|-----------------------------|--|---------------|-------------------------|-------|------------------------|-------|----------|---|-----------------------|-----------------------|---------|
| 108-88-3 | Toluene | 5.20 | | ug/L | 0.346 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PA | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 156-60-5 | trans-1,2-Dichloroethylene | 144 | | ug/L | 55.8 | 100 | 200 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PA | 03/21/2023 06:13 | 03/22/2023 02:23 | JTG |
| 10061-02-6 | trans-1,3-Dichloropropylene | ND | | ug/L | 0.229 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PA | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 79-01-6 | Trichloroethylene | 196 | | ug/L | 49.8 | 100 | 200 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PA | 03/21/2023 06:13 | 03/22/2023 02:23 | JTG |
| 75-69-4 | Trichlorofluoromethane | ND | | ug/L | 0.337 | 0.500 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PA | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| 75-01-4 | Vinyl Chloride | 4640 | | ug/L | 93.8 | 100 | 200 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP,PA | 03/21/2023 06:13 | 03/22/2023 02:23 | JTG |
| 1330-20-7 | Xylenes, Total | 2.72 | | ug/L | 0.836 | 1.50 | 1 | EPA 8260C Certifications: CTDOH-PH-0723,NELAC-NY10854,NELAC-NY12058,NJDEP | 03/20/2023 06:45 | 03/21/2023 09:54 | JTG |
| Surrogate Recoveries | | Result | Acceptance Range | | | | | | | | |
| 17060-07-0 | Surrogate: SURRE: 1,2-Dichloroethane-d4 | 104 % | 69-130 | | | | | | | | |
| 2037-26-5 | Surrogate: SURRE: Toluene-d8 | 101 % | 81-117 | | | | | | | | |
| 460-00-4 | Surrogate: SURRE: p-Bromofluorobenzene | 92.0 % | 79-122 | | | | | | | | |



Analytical Batch Summary

Batch ID: BC31334

Preparation Method: EPA 5030B

Prepared By: JTG

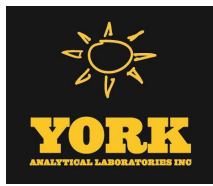
| YORK Sample ID | Client Sample ID | Preparation Date |
|----------------|------------------|------------------|
| 23C0637-01 | DW-1 | 03/20/23 |
| 23C0637-02 | DW-2 | 03/20/23 |
| BC31334-BLK1 | Blank | 03/20/23 |
| BC31334-BS1 | LCS | 03/20/23 |
| BC31334-BSD1 | LCS Dup | 03/20/23 |

Batch ID: BC31487

Preparation Method: EPA 5030B

Prepared By: JTG

| YORK Sample ID | Client Sample ID | Preparation Date |
|----------------|------------------|------------------|
| 23C0637-02RE1 | DW-2 | 03/21/23 |
| BC31487-BLK1 | Blank | 03/21/23 |
| BC31487-BS1 | LCS | 03/21/23 |
| BC31487-BSD1 | LCS Dup | 03/21/23 |



Volatile Organic Compounds by GC/MS - Quality Control Data
York Analytical Laboratories, Inc. - Stratford

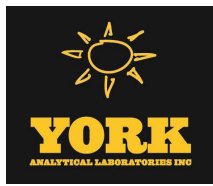
| Analyte | Result | Reporting Limit | Units | Spike Level | Source* Result | %REC | %REC Limits | Flag | RPD | RPD Limit | Flag |
|---------|--------|-----------------|-------|-------------|----------------|------|-------------|------|-----|-----------|------|
|---------|--------|-----------------|-------|-------------|----------------|------|-------------|------|-----|-----------|------|

Batch BC31334 - EPA 5030B

Blank (BC31334-BLK1)

Prepared: 03/20/2023 Analyzed: 03/21/2023

| | | | | | | | | | | | |
|---|----|-------|------|--|--|--|--|--|--|--|--|
| 1,1,1,2-Tetrachloroethane | ND | 0.500 | ug/L | | | | | | | | |
| 1,1,1-Trichloroethane | ND | 0.500 | " | | | | | | | | |
| 1,1,2,2-Tetrachloroethane | ND | 0.500 | " | | | | | | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | ND | 0.500 | " | | | | | | | | |
| 1,1,2-Trichloroethane | ND | 0.500 | " | | | | | | | | |
| 1,1-Dichloroethane | ND | 0.500 | " | | | | | | | | |
| 1,1-Dichloroethylene | ND | 0.500 | " | | | | | | | | |
| 1,2,3-Trichlorobenzene | ND | 0.500 | " | | | | | | | | |
| 1,2,3-Trichloropropane | ND | 0.500 | " | | | | | | | | |
| 1,2,4-Trichlorobenzene | ND | 0.500 | " | | | | | | | | |
| 1,2,4-Trimethylbenzene | ND | 0.500 | " | | | | | | | | |
| 1,2-Dibromo-3-chloropropane | ND | 0.500 | " | | | | | | | | |
| 1,2-Dibromoethane | ND | 0.500 | " | | | | | | | | |
| 1,2-Dichlorobenzene | ND | 0.500 | " | | | | | | | | |
| 1,2-Dichloroethane | ND | 0.500 | " | | | | | | | | |
| 1,2-Dichloropropane | ND | 0.500 | " | | | | | | | | |
| 1,3,5-Trimethylbenzene | ND | 0.500 | " | | | | | | | | |
| 1,3-Dichlorobenzene | ND | 0.500 | " | | | | | | | | |
| 1,4-Dichlorobenzene | ND | 0.500 | " | | | | | | | | |
| 1,4-Dioxane | ND | 80.0 | " | | | | | | | | |
| 2-Butanone | ND | 0.500 | " | | | | | | | | |
| 2-Hexanone | ND | 0.500 | " | | | | | | | | |
| 4-Methyl-2-pentanone | ND | 0.500 | " | | | | | | | | |
| Acetone | ND | 2.00 | " | | | | | | | | |
| Acrolein | ND | 0.500 | " | | | | | | | | |
| Acrylonitrile | ND | 0.500 | " | | | | | | | | |
| Benzene | ND | 0.500 | " | | | | | | | | |
| Bromochloromethane | ND | 0.500 | " | | | | | | | | |
| Bromodichloromethane | ND | 0.500 | " | | | | | | | | |
| Bromoform | ND | 0.500 | " | | | | | | | | |
| Bromomethane | ND | 0.500 | " | | | | | | | | |
| Carbon disulfide | ND | 0.500 | " | | | | | | | | |
| Carbon tetrachloride | ND | 0.500 | " | | | | | | | | |
| Chlorobenzene | ND | 0.500 | " | | | | | | | | |
| Chloroethane | ND | 0.500 | " | | | | | | | | |
| Chloroform | ND | 0.500 | " | | | | | | | | |
| Chloromethane | ND | 0.500 | " | | | | | | | | |
| cis-1,2-Dichloroethylene | ND | 0.500 | " | | | | | | | | |
| cis-1,3-Dichloropropylene | ND | 0.500 | " | | | | | | | | |
| Cyclohexane | ND | 0.500 | " | | | | | | | | |
| Dibromochloromethane | ND | 0.500 | " | | | | | | | | |
| Dibromomethane | ND | 0.500 | " | | | | | | | | |
| Dichlorodifluoromethane | ND | 0.500 | " | | | | | | | | |
| Ethyl Benzene | ND | 0.500 | " | | | | | | | | |
| Hexachlorobutadiene | ND | 0.500 | " | | | | | | | | |
| Isopropylbenzene | ND | 0.500 | " | | | | | | | | |
| Methyl acetate | ND | 0.500 | " | | | | | | | | |
| Methyl tert-butyl ether (MTBE) | ND | 0.500 | " | | | | | | | | |
| Methylcyclohexane | ND | 0.500 | " | | | | | | | | |



Volatile Organic Compounds by GC/MS - Quality Control Data
York Analytical Laboratories, Inc. - Stratford

| Analyte | Result | Reporting Limit | Units | Spike Level | Source* Result | %REC | %REC Limits | Flag | RPD | RPD Limit | Flag |
|---------|--------|-----------------|-------|-------------|----------------|------|-------------|------|-----|-----------|------|
|---------|--------|-----------------|-------|-------------|----------------|------|-------------|------|-----|-----------|------|

Batch BC31334 - EPA 5030B

Blank (BC31334-BLK1)

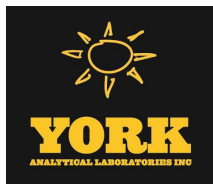
Prepared: 03/20/2023 Analyzed: 03/21/2023

| | | | | | | | | | | | |
|--|------|-------|------|------|--|------|--------|--|--|--|--|
| Methylene chloride | ND | 2.00 | ug/L | | | | | | | | |
| n-Butylbenzene | ND | 0.500 | " | | | | | | | | |
| n-Propylbenzene | ND | 0.500 | " | | | | | | | | |
| o-Xylene | ND | 0.500 | " | | | | | | | | |
| p- & m- Xylenes | ND | 1.00 | " | | | | | | | | |
| p-Isopropyltoluene | ND | 0.500 | " | | | | | | | | |
| sec-Butylbenzene | ND | 0.500 | " | | | | | | | | |
| Styrene | ND | 0.500 | " | | | | | | | | |
| tert-Butyl alcohol (TBA) | ND | 1.00 | " | | | | | | | | |
| tert-Butylbenzene | ND | 0.500 | " | | | | | | | | |
| Tetrachloroethylene | ND | 0.500 | " | | | | | | | | |
| Toluene | ND | 0.500 | " | | | | | | | | |
| trans-1,2-Dichloroethylene | ND | 0.500 | " | | | | | | | | |
| trans-1,3-Dichloropropylene | ND | 0.500 | " | | | | | | | | |
| Trichloroethylene | ND | 0.500 | " | | | | | | | | |
| Trichlorofluoromethane | ND | 0.500 | " | | | | | | | | |
| Vinyl Chloride | ND | 0.500 | " | | | | | | | | |
| Xylenes, Total | ND | 1.50 | " | | | | | | | | |
| <i>Surrogate: SURRE: 1,2-Dichloroethane-d4</i> | 9.80 | | " | 10.0 | | 98.0 | 69-130 | | | | |
| <i>Surrogate: SURRE: Toluene-d8</i> | 10.0 | | " | 10.0 | | 100 | 81-117 | | | | |
| <i>Surrogate: SURRE: p-Bromofluorobenzene</i> | 10.3 | | " | 10.0 | | 103 | 79-122 | | | | |

LCS (BC31334-BS1)

Prepared & Analyzed: 03/20/2023

| | | | | | | | | | | | |
|---|------|--|------|------|--|------|--------|--|--|--|--|
| 1,1,1,2-Tetrachloroethane | 9.85 | | ug/L | 10.0 | | 98.5 | 82-126 | | | | |
| 1,1,1-Trichloroethane | 10.0 | | " | 10.0 | | 100 | 78-136 | | | | |
| 1,1,2,2-Tetrachloroethane | 10.8 | | " | 10.0 | | 108 | 76-129 | | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 11.6 | | " | 10.0 | | 116 | 54-165 | | | | |
| 1,1,2-Trichloroethane | 10.2 | | " | 10.0 | | 102 | 82-123 | | | | |
| 1,1-Dichloroethane | 10.9 | | " | 10.0 | | 109 | 82-129 | | | | |
| 1,1-Dichloroethylene | 11.6 | | " | 10.0 | | 116 | 68-138 | | | | |
| 1,2,3-Trichlorobenzene | 8.19 | | " | 10.0 | | 81.9 | 76-136 | | | | |
| 1,2,3-Trichloropropane | 10.1 | | " | 10.0 | | 101 | 77-128 | | | | |
| 1,2,4-Trichlorobenzene | 8.51 | | " | 10.0 | | 85.1 | 76-137 | | | | |
| 1,2,4-Trimethylbenzene | 10.9 | | " | 10.0 | | 109 | 82-132 | | | | |
| 1,2-Dibromo-3-chloropropane | 9.45 | | " | 10.0 | | 94.5 | 45-147 | | | | |
| 1,2-Dibromoethane | 9.80 | | " | 10.0 | | 98.0 | 83-124 | | | | |
| 1,2-Dichlorobenzene | 10.3 | | " | 10.0 | | 103 | 79-123 | | | | |
| 1,2-Dichloroethane | 10.3 | | " | 10.0 | | 103 | 73-132 | | | | |
| 1,2-Dichloropropane | 12.3 | | " | 10.0 | | 123 | 78-126 | | | | |
| 1,3,5-Trimethylbenzene | 11.2 | | " | 10.0 | | 112 | 80-131 | | | | |
| 1,3-Dichlorobenzene | 10.3 | | " | 10.0 | | 103 | 86-122 | | | | |
| 1,4-Dichlorobenzene | 10.1 | | " | 10.0 | | 101 | 85-124 | | | | |
| 1,4-Dioxane | 248 | | " | 210 | | 118 | 10-349 | | | | |
| 2-Butanone | 10.0 | | " | 10.0 | | 100 | 49-152 | | | | |
| 2-Hexanone | 9.46 | | " | 10.0 | | 94.6 | 51-146 | | | | |
| 4-Methyl-2-pentanone | 8.77 | | " | 10.0 | | 87.7 | 57-145 | | | | |
| Acetone | 10.6 | | " | 10.0 | | 106 | 14-150 | | | | |
| Acrolein | 6.55 | | " | 10.0 | | 65.5 | 10-153 | | | | |
| Acrylonitrile | 9.49 | | " | 10.0 | | 94.9 | 51-150 | | | | |
| Benzene | 11.7 | | " | 10.0 | | 117 | 85-126 | | | | |



Volatile Organic Compounds by GC/MS - Quality Control Data
York Analytical Laboratories, Inc. - Stratford

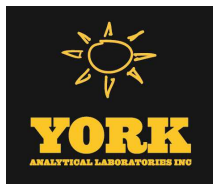
| Analyte | Result | Reporting Limit | Units | Spike Level | Source* Result | %REC | %REC Limits | Flag | RPD | RPD Limit | Flag |
|---------|--------|-----------------|-------|-------------|----------------|------|-------------|------|-----|-----------|------|
|---------|--------|-----------------|-------|-------------|----------------|------|-------------|------|-----|-----------|------|

Batch BC31334 - EPA 5030B

LCS (BC31334-BS1)

Prepared & Analyzed: 03/20/2023

| | | | | | | | | | | | |
|---|------|--|------|------|--|------|--------|----------|--|--|--|
| Bromochloromethane | 11.2 | | ug/L | 10.0 | | 112 | 77-128 | | | | |
| Bromodichloromethane | 10.4 | | " | 10.0 | | 104 | 79-128 | | | | |
| Bromoform | 8.05 | | " | 10.0 | | 80.5 | 78-133 | | | | |
| Bromomethane | 8.87 | | " | 10.0 | | 88.7 | 43-168 | | | | |
| Carbon disulfide | 11.4 | | " | 10.0 | | 114 | 68-146 | | | | |
| Carbon tetrachloride | 10.4 | | " | 10.0 | | 104 | 77-141 | | | | |
| Chlorobenzene | 11.3 | | " | 10.0 | | 113 | 88-120 | | | | |
| Chloroethane | 9.52 | | " | 10.0 | | 95.2 | 65-136 | | | | |
| Chloroform | 10.6 | | " | 10.0 | | 106 | 82-128 | | | | |
| Chloromethane | 10.5 | | " | 10.0 | | 105 | 43-155 | | | | |
| cis-1,2-Dichloroethylene | 11.0 | | " | 10.0 | | 110 | 83-129 | | | | |
| cis-1,3-Dichloropropylene | 10.5 | | " | 10.0 | | 105 | 80-131 | | | | |
| Cyclohexane | 5.76 | | " | 10.0 | | 57.6 | 63-149 | Low Bias | | | |
| Dibromochloromethane | 9.10 | | " | 10.0 | | 91.0 | 80-130 | | | | |
| Dibromomethane | 10.2 | | " | 10.0 | | 102 | 72-134 | | | | |
| Dichlorodifluoromethane | 9.73 | | " | 10.0 | | 97.3 | 44-144 | | | | |
| Ethyl Benzene | 11.9 | | " | 10.0 | | 119 | 80-131 | | | | |
| Hexachlorobutadiene | 7.39 | | " | 10.0 | | 73.9 | 67-146 | | | | |
| Isopropylbenzene | 11.7 | | " | 10.0 | | 117 | 76-140 | | | | |
| Methyl acetate | 10.2 | | " | 10.0 | | 102 | 51-139 | | | | |
| Methyl tert-butyl ether (MTBE) | 9.13 | | " | 10.0 | | 91.3 | 76-135 | | | | |
| Methylcyclohexane | 12.1 | | " | 10.0 | | 121 | 72-143 | | | | |
| Methylene chloride | 11.1 | | " | 10.0 | | 111 | 55-137 | | | | |
| n-Butylbenzene | 11.6 | | " | 10.0 | | 116 | 79-132 | | | | |
| n-Propylbenzene | 12.3 | | " | 10.0 | | 123 | 78-133 | | | | |
| o-Xylene | 11.5 | | " | 10.0 | | 115 | 78-130 | | | | |
| p- & m- Xylenes | 23.5 | | " | 20.0 | | 118 | 77-133 | | | | |
| p-Isopropyltoluene | 11.1 | | " | 10.0 | | 111 | 81-136 | | | | |
| sec-Butylbenzene | 11.5 | | " | 10.0 | | 115 | 79-137 | | | | |
| Styrene | 11.2 | | " | 10.0 | | 112 | 67-132 | | | | |
| tert-Butyl alcohol (TBA) | 24.5 | | " | 50.0 | | 49.1 | 25-162 | | | | |
| tert-Butylbenzene | 9.68 | | " | 10.0 | | 96.8 | 77-138 | | | | |
| Tetrachloroethylene | 10.2 | | " | 10.0 | | 102 | 82-131 | | | | |
| Toluene | 11.3 | | " | 10.0 | | 113 | 80-127 | | | | |
| trans-1,2-Dichloroethylene | 11.3 | | " | 10.0 | | 113 | 80-132 | | | | |
| trans-1,3-Dichloropropylene | 9.81 | | " | 10.0 | | 98.1 | 78-131 | | | | |
| Trichloroethylene | 11.1 | | " | 10.0 | | 111 | 82-128 | | | | |
| Trichlorofluoromethane | 12.3 | | " | 10.0 | | 123 | 67-139 | | | | |
| Vinyl Chloride | 9.74 | | " | 10.0 | | 97.4 | 58-145 | | | | |
| Surrogate: SURRE: 1,2-Dichloroethane-d4 | 9.46 | | " | 10.0 | | 94.6 | 69-130 | | | | |
| Surrogate: SURRE: Toluene-d8 | 10.0 | | " | 10.0 | | 100 | 81-117 | | | | |
| Surrogate: SURRE: p-Bromofluorobenzene | 10.2 | | " | 10.0 | | 102 | 79-122 | | | | |



Volatile Organic Compounds by GC/MS - Quality Control Data
York Analytical Laboratories, Inc. - Stratford

| Analyte | Result | Reporting Limit | Units | Spike Level | Source* Result | %REC | %REC Limits | Flag | RPD | RPD Limit | Flag |
|---|--------|-----------------|-------|-------------|----------------|------|-------------|----------|-------|---|------|
| Batch BC31334 - EPA 5030B | | | | | | | | | | | |
| LCS Dup (BC31334-BSD1) | | | | | | | | | | | |
| | | | | | | | | | | Prepared: 03/20/2023 Analyzed: 03/21/2023 | |
| 1,1,1,2-Tetrachloroethane | 9.64 | | ug/L | 10.0 | | 96.4 | 82-126 | | 2.15 | 30 | |
| 1,1,1-Trichloroethane | 9.73 | | " | 10.0 | | 97.3 | 78-136 | | 2.74 | 30 | |
| 1,1,2,2-Tetrachloroethane | 10.6 | | " | 10.0 | | 106 | 76-129 | | 1.22 | 30 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 11.1 | | " | 10.0 | | 111 | 54-165 | | 4.76 | 30 | |
| 1,1,2-Trichloroethane | 10.2 | | " | 10.0 | | 102 | 82-123 | | 0.293 | 30 | |
| 1,1-Dichloroethane | 10.6 | | " | 10.0 | | 106 | 82-129 | | 2.69 | 30 | |
| 1,1-Dichloroethylene | 11.0 | | " | 10.0 | | 110 | 68-138 | | 5.92 | 30 | |
| 1,2,3-Trichlorobenzene | 8.03 | | " | 10.0 | | 80.3 | 76-136 | | 1.97 | 30 | |
| 1,2,3-Trichloropropane | 9.95 | | " | 10.0 | | 99.5 | 77-128 | | 1.40 | 30 | |
| 1,2,4-Trichlorobenzene | 8.42 | | " | 10.0 | | 84.2 | 76-137 | | 1.06 | 30 | |
| 1,2,4-Trimethylbenzene | 10.3 | | " | 10.0 | | 103 | 82-132 | | 5.37 | 30 | |
| 1,2-Dibromo-3-chloropropane | 9.24 | | " | 10.0 | | 92.4 | 45-147 | | 2.25 | 30 | |
| 1,2-Dibromoethane | 9.91 | | " | 10.0 | | 99.1 | 83-124 | | 1.12 | 30 | |
| 1,2-Dichlorobenzene | 9.92 | | " | 10.0 | | 99.2 | 79-123 | | 3.86 | 30 | |
| 1,2-Dichloroethane | 10.3 | | " | 10.0 | | 103 | 73-132 | | 0.582 | 30 | |
| 1,2-Dichloropropane | 12.0 | | " | 10.0 | | 120 | 78-126 | | 1.98 | 30 | |
| 1,3,5-Trimethylbenzene | 10.6 | | " | 10.0 | | 106 | 80-131 | | 6.33 | 30 | |
| 1,3-Dichlorobenzene | 9.72 | | " | 10.0 | | 97.2 | 86-122 | | 5.70 | 30 | |
| 1,4-Dichlorobenzene | 9.73 | | " | 10.0 | | 97.3 | 85-124 | | 3.34 | 30 | |
| 1,4-Dioxane | 242 | | " | 210 | | 115 | 10-349 | | 2.73 | 30 | |
| 2-Butanone | 10.3 | | " | 10.0 | | 103 | 49-152 | | 2.84 | 30 | |
| 2-Hexanone | 9.60 | | " | 10.0 | | 96.0 | 51-146 | | 1.47 | 30 | |
| 4-Methyl-2-pentanone | 8.80 | | " | 10.0 | | 88.0 | 57-145 | | 0.341 | 30 | |
| Acetone | 10.7 | | " | 10.0 | | 107 | 14-150 | | 0.845 | 30 | |
| Acrolein | 6.73 | | " | 10.0 | | 67.3 | 10-153 | | 2.71 | 30 | |
| Acrylonitrile | 9.95 | | " | 10.0 | | 99.5 | 51-150 | | 4.73 | 30 | |
| Benzene | 11.3 | | " | 10.0 | | 113 | 85-126 | | 2.87 | 30 | |
| Bromochloromethane | 11.2 | | " | 10.0 | | 112 | 77-128 | | 0.00 | 30 | |
| Bromodichloromethane | 10.2 | | " | 10.0 | | 102 | 79-128 | | 1.84 | 30 | |
| Bromoform | 8.13 | | " | 10.0 | | 81.3 | 78-133 | | 0.989 | 30 | |
| Bromomethane | 8.86 | | " | 10.0 | | 88.6 | 43-168 | | 0.113 | 30 | |
| Carbon disulfide | 10.8 | | " | 10.0 | | 108 | 68-146 | | 5.41 | 30 | |
| Carbon tetrachloride | 9.83 | | " | 10.0 | | 98.3 | 77-141 | | 5.15 | 30 | |
| Chlorobenzene | 10.9 | | " | 10.0 | | 109 | 88-120 | | 3.61 | 30 | |
| Chloroethane | 8.94 | | " | 10.0 | | 89.4 | 65-136 | | 6.28 | 30 | |
| Chloroform | 10.3 | | " | 10.0 | | 103 | 82-128 | | 2.49 | 30 | |
| Chloromethane | 9.94 | | " | 10.0 | | 99.4 | 43-155 | | 5.57 | 30 | |
| cis-1,2-Dichloroethylene | 10.7 | | " | 10.0 | | 107 | 83-129 | | 2.86 | 30 | |
| cis-1,3-Dichloropropylene | 10.3 | | " | 10.0 | | 103 | 80-131 | | 1.83 | 30 | |
| Cyclohexane | 5.44 | | " | 10.0 | | 54.4 | 63-149 | Low Bias | 5.71 | 30 | |
| Dibromochloromethane | 9.14 | | " | 10.0 | | 91.4 | 80-130 | | 0.439 | 30 | |
| Dibromomethane | 10.4 | | " | 10.0 | | 104 | 72-134 | | 2.24 | 30 | |
| Dichlorodifluoromethane | 9.04 | | " | 10.0 | | 90.4 | 44-144 | | 7.35 | 30 | |
| Ethyl Benzene | 11.4 | | " | 10.0 | | 114 | 80-131 | | 4.48 | 30 | |
| Hexachlorobutadiene | 7.27 | | " | 10.0 | | 72.7 | 67-146 | | 1.64 | 30 | |
| Isopropylbenzene | 10.9 | | " | 10.0 | | 109 | 76-140 | | 6.80 | 30 | |
| Methyl acetate | 10.7 | | " | 10.0 | | 107 | 51-139 | | 4.86 | 30 | |
| Methyl tert-butyl ether (MTBE) | 9.46 | | " | 10.0 | | 94.6 | 76-135 | | 3.55 | 30 | |
| Methylcyclohexane | 11.3 | | " | 10.0 | | 113 | 72-143 | | 6.50 | 30 | |
| Methylene chloride | 11.0 | | " | 10.0 | | 110 | 55-137 | | 0.811 | 30 | |



Volatile Organic Compounds by GC/MS - Quality Control Data
York Analytical Laboratories, Inc. - Stratford

| Analyte | Result | Reporting Limit | Units | Spike Level | Source* Result | %REC | %REC Limits | Flag | RPD | RPD Limit | Flag |
|---------|--------|-----------------|-------|-------------|----------------|------|-------------|------|-----|-----------|------|
|---------|--------|-----------------|-------|-------------|----------------|------|-------------|------|-----|-----------|------|

Batch BC31334 - EPA 5030B

LCS Dup (BC31334-BSD1)

Prepared: 03/20/2023 Analyzed: 03/21/2023

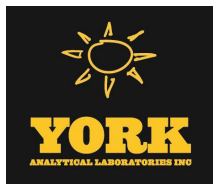
| | | | | | | | | | | | |
|---|-------------|--|----------|-------------|--|-------------|---------------|--|-------|----|--|
| n-Butylbenzene | 10.8 | | ug/L | 10.0 | | 108 | 79-132 | | 7.03 | 30 | |
| n-Propylbenzene | 11.5 | | " | 10.0 | | 115 | 78-133 | | 6.65 | 30 | |
| o-Xylene | 11.1 | | " | 10.0 | | 111 | 78-130 | | 3.63 | 30 | |
| p- & m- Xylenes | 22.8 | | " | 20.0 | | 114 | 77-133 | | 3.41 | 30 | |
| p-Isopropyltoluene | 10.3 | | " | 10.0 | | 103 | 81-136 | | 7.96 | 30 | |
| sec-Butylbenzene | 10.8 | | " | 10.0 | | 108 | 79-137 | | 6.74 | 30 | |
| Styrene | 10.9 | | " | 10.0 | | 109 | 67-132 | | 3.26 | 30 | |
| tert-Butyl alcohol (TBA) | 26.6 | | " | 50.0 | | 53.3 | 25-162 | | 8.21 | 30 | |
| tert-Butylbenzene | 9.03 | | " | 10.0 | | 90.3 | 77-138 | | 6.95 | 30 | |
| Tetrachloroethylene | 9.63 | | " | 10.0 | | 96.3 | 82-131 | | 5.36 | 30 | |
| Toluene | 10.8 | | " | 10.0 | | 108 | 80-127 | | 4.34 | 30 | |
| trans-1,2-Dichloroethylene | 10.8 | | " | 10.0 | | 108 | 80-132 | | 4.52 | 30 | |
| trans-1,3-Dichloropropylene | 9.80 | | " | 10.0 | | 98.0 | 78-131 | | 0.102 | 30 | |
| Trichloroethylene | 10.6 | | " | 10.0 | | 106 | 82-128 | | 5.17 | 30 | |
| Trichlorofluoromethane | 11.7 | | " | 10.0 | | 117 | 67-139 | | 5.33 | 30 | |
| Vinyl Chloride | 9.24 | | " | 10.0 | | 92.4 | 58-145 | | 5.27 | 30 | |
| <i>Surrogate: SURR: 1,2-Dichloroethane-d4</i> | <i>9.77</i> | | <i>"</i> | <i>10.0</i> | | <i>97.7</i> | <i>69-130</i> | | | | |
| <i>Surrogate: SURR: Toluene-d8</i> | <i>9.96</i> | | <i>"</i> | <i>10.0</i> | | <i>99.6</i> | <i>81-117</i> | | | | |
| <i>Surrogate: SURR: p-Bromofluorobenzene</i> | <i>10.1</i> | | <i>"</i> | <i>10.0</i> | | <i>101</i> | <i>79-122</i> | | | | |

Batch BC31487 - EPA 5030B

Blank (BC31487-BLK1)

Prepared & Analyzed: 03/21/2023

| | | | | | | | | | | | |
|---|----|-------|------|--|--|--|--|--|--|--|--|
| 1,1,1,2-Tetrachloroethane | ND | 0.500 | ug/L | | | | | | | | |
| 1,1,1-Trichloroethane | ND | 0.500 | " | | | | | | | | |
| 1,1,2,2-Tetrachloroethane | ND | 0.500 | " | | | | | | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | ND | 0.500 | " | | | | | | | | |
| 1,1,2-Trichloroethane | ND | 0.500 | " | | | | | | | | |
| 1,1-Dichloroethane | ND | 0.500 | " | | | | | | | | |
| 1,1-Dichloroethylene | ND | 0.500 | " | | | | | | | | |
| 1,2,3-Trichlorobenzene | ND | 0.500 | " | | | | | | | | |
| 1,2,3-Trichloropropane | ND | 0.500 | " | | | | | | | | |
| 1,2,4-Trichlorobenzene | ND | 0.500 | " | | | | | | | | |
| 1,2,4-Trimethylbenzene | ND | 0.500 | " | | | | | | | | |
| 1,2-Dibromo-3-chloropropane | ND | 0.500 | " | | | | | | | | |
| 1,2-Dibromoethane | ND | 0.500 | " | | | | | | | | |
| 1,2-Dichlorobenzene | ND | 0.500 | " | | | | | | | | |
| 1,2-Dichloroethane | ND | 0.500 | " | | | | | | | | |
| 1,2-Dichloropropane | ND | 0.500 | " | | | | | | | | |
| 1,3,5-Trimethylbenzene | ND | 0.500 | " | | | | | | | | |
| 1,3-Dichlorobenzene | ND | 0.500 | " | | | | | | | | |
| 1,4-Dichlorobenzene | ND | 0.500 | " | | | | | | | | |
| 1,4-Dioxane | ND | 80.0 | " | | | | | | | | |
| 2-Butanone | ND | 0.500 | " | | | | | | | | |
| 2-Hexanone | ND | 0.500 | " | | | | | | | | |
| 4-Methyl-2-pentanone | ND | 0.500 | " | | | | | | | | |
| Acetone | ND | 2.00 | " | | | | | | | | |
| Acrolein | ND | 0.500 | " | | | | | | | | |
| Acrylonitrile | ND | 0.500 | " | | | | | | | | |
| Benzene | ND | 0.500 | " | | | | | | | | |
| Bromochloromethane | ND | 0.500 | " | | | | | | | | |



Volatile Organic Compounds by GC/MS - Quality Control Data
York Analytical Laboratories, Inc. - Stratford

| Analyte | Result | Reporting | Units | Spike Level | Source* Result | %REC | %REC Limits | Flag | RPD | RPD | |
|---------|--------|-----------|-------|----------------|-------------------|------|----------------|------|-----|-------|------|
| | | Limit | | | | | | | | Limit | Flag |

Batch BC31487 - EPA 5030B

Blank (BC31487-BLK1)

Prepared & Analyzed: 03/21/2023

| | | | | | | | | | | | |
|---|------|-------|------|------|--|------|--------|--|--|--|--|
| Bromodichloromethane | ND | 0.500 | ug/L | | | | | | | | |
| Bromoform | ND | 0.500 | " | | | | | | | | |
| Bromomethane | ND | 0.500 | " | | | | | | | | |
| Carbon disulfide | ND | 0.500 | " | | | | | | | | |
| Carbon tetrachloride | ND | 0.500 | " | | | | | | | | |
| Chlorobenzene | ND | 0.500 | " | | | | | | | | |
| Chloroethane | ND | 0.500 | " | | | | | | | | |
| Chloroform | ND | 0.500 | " | | | | | | | | |
| Chloromethane | ND | 0.500 | " | | | | | | | | |
| cis-1,2-Dichloroethylene | ND | 0.500 | " | | | | | | | | |
| cis-1,3-Dichloropropylene | ND | 0.500 | " | | | | | | | | |
| Cyclohexane | ND | 0.500 | " | | | | | | | | |
| Dibromochloromethane | ND | 0.500 | " | | | | | | | | |
| Dibromomethane | ND | 0.500 | " | | | | | | | | |
| Dichlorodifluoromethane | ND | 0.500 | " | | | | | | | | |
| Ethyl Benzene | ND | 0.500 | " | | | | | | | | |
| Hexachlorobutadiene | ND | 0.500 | " | | | | | | | | |
| Isopropylbenzene | ND | 0.500 | " | | | | | | | | |
| Methyl acetate | ND | 0.500 | " | | | | | | | | |
| Methyl tert-butyl ether (MTBE) | ND | 0.500 | " | | | | | | | | |
| Methylcyclohexane | ND | 0.500 | " | | | | | | | | |
| Methylene chloride | ND | 2.00 | " | | | | | | | | |
| n-Butylbenzene | ND | 0.500 | " | | | | | | | | |
| n-Propylbenzene | ND | 0.500 | " | | | | | | | | |
| o-Xylene | ND | 0.500 | " | | | | | | | | |
| p- & m- Xylenes | ND | 1.00 | " | | | | | | | | |
| p-Isopropyltoluene | ND | 0.500 | " | | | | | | | | |
| sec-Butylbenzene | ND | 0.500 | " | | | | | | | | |
| Styrene | ND | 0.500 | " | | | | | | | | |
| tert-Butyl alcohol (TBA) | ND | 1.00 | " | | | | | | | | |
| tert-Butylbenzene | ND | 0.500 | " | | | | | | | | |
| Tetrachloroethylene | ND | 0.500 | " | | | | | | | | |
| Toluene | ND | 0.500 | " | | | | | | | | |
| trans-1,2-Dichloroethylene | ND | 0.500 | " | | | | | | | | |
| trans-1,3-Dichloropropylene | ND | 0.500 | " | | | | | | | | |
| Trichloroethylene | ND | 0.500 | " | | | | | | | | |
| Trichlorofluoromethane | ND | 0.500 | " | | | | | | | | |
| Vinyl Chloride | ND | 0.500 | " | | | | | | | | |
| Xylenes, Total | ND | 1.50 | " | | | | | | | | |
| <hr/> | | | | | | | | | | | |
| Surrogate: SURRE: 1,2-Dichloroethane-d4 | 9.81 | | " | 10.0 | | 98.1 | 69-130 | | | | |
| Surrogate: SURRE: Toluene-d8 | 10.9 | | " | 10.0 | | 109 | 81-117 | | | | |
| Surrogate: SURRE: p-Bromofluorobenzene | 9.60 | | " | 10.0 | | 96.0 | 79-122 | | | | |



Volatile Organic Compounds by GC/MS - Quality Control Data

York Analytical Laboratories, Inc. - Stratford

| Analyte | Result | Reporting Limit | Units | Spike Level | Source* Result | %REC | %REC Limits | Flag | RPD | RPD Limit | Flag |
|---|--------|-----------------|-------|-------------|----------------|------|-------------|-----------|-----|-----------|------|
| Batch BC31487 - EPA 5030B | | | | | | | | | | | |
| LCS (BC31487-BS1) | | | | | | | | | | | |
| Prepared & Analyzed: 03/21/2023 | | | | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | 10.1 | | ug/L | 10.0 | | 101 | 82-126 | | | | |
| 1,1,1-Trichloroethane | 9.04 | | " | 10.0 | | 90.4 | 78-136 | | | | |
| 1,1,2,2-Tetrachloroethane | 11.1 | | " | 10.0 | | 111 | 76-129 | | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 10.4 | | " | 10.0 | | 104 | 54-165 | | | | |
| 1,1,2-Trichloroethane | 10.4 | | " | 10.0 | | 104 | 82-123 | | | | |
| 1,1-Dichloroethane | 10.0 | | " | 10.0 | | 100 | 82-129 | | | | |
| 1,1-Dichloroethylene | 10.6 | | " | 10.0 | | 106 | 68-138 | | | | |
| 1,2,3-Trichlorobenzene | 11.1 | | " | 10.0 | | 111 | 76-136 | | | | |
| 1,2,3-Trichloropropane | 10.0 | | " | 10.0 | | 100 | 77-128 | | | | |
| 1,2,4-Trichlorobenzene | 10.3 | | " | 10.0 | | 103 | 76-137 | | | | |
| 1,2,4-Trimethylbenzene | 11.4 | | " | 10.0 | | 114 | 82-132 | | | | |
| 1,2-Dibromo-3-chloropropane | 10.9 | | " | 10.0 | | 109 | 45-147 | | | | |
| 1,2-Dibromoethane | 9.68 | | " | 10.0 | | 96.8 | 83-124 | | | | |
| 1,2-Dichlorobenzene | 10.9 | | " | 10.0 | | 109 | 79-123 | | | | |
| 1,2-Dichloroethane | 9.06 | | " | 10.0 | | 90.6 | 73-132 | | | | |
| 1,2-Dichloropropane | 12.5 | | " | 10.0 | | 125 | 78-126 | | | | |
| 1,3,5-Trimethylbenzene | 11.7 | | " | 10.0 | | 117 | 80-131 | | | | |
| 1,3-Dichlorobenzene | 10.2 | | " | 10.0 | | 102 | 86-122 | | | | |
| 1,4-Dichlorobenzene | 10.0 | | " | 10.0 | | 100 | 85-124 | | | | |
| 1,4-Dioxane | 197 | | " | 210 | | 93.9 | 10-349 | | | | |
| 2-Butanone | 9.64 | | " | 10.0 | | 96.4 | 49-152 | | | | |
| 2-Hexanone | 9.42 | | " | 10.0 | | 94.2 | 51-146 | | | | |
| 4-Methyl-2-pentanone | 8.78 | | " | 10.0 | | 87.8 | 57-145 | | | | |
| Acetone | 9.66 | | " | 10.0 | | 96.6 | 14-150 | | | | |
| Acrolein | 5.49 | | " | 10.0 | | 54.9 | 10-153 | | | | |
| Acrylonitrile | 8.52 | | " | 10.0 | | 85.2 | 51-150 | | | | |
| Benzene | 10.7 | | " | 10.0 | | 107 | 85-126 | | | | |
| Bromochloromethane | 10.1 | | " | 10.0 | | 101 | 77-128 | | | | |
| Bromodichloromethane | 10.6 | | " | 10.0 | | 106 | 79-128 | | | | |
| Bromoform | 8.24 | | " | 10.0 | | 82.4 | 78-133 | | | | |
| Bromomethane | 8.96 | | " | 10.0 | | 89.6 | 43-168 | | | | |
| Carbon disulfide | 10.4 | | " | 10.0 | | 104 | 68-146 | | | | |
| Carbon tetrachloride | 9.23 | | " | 10.0 | | 92.3 | 77-141 | | | | |
| Chlorobenzene | 11.4 | | " | 10.0 | | 114 | 88-120 | | | | |
| Chloroethane | 9.07 | | " | 10.0 | | 90.7 | 65-136 | | | | |
| Chloroform | 9.56 | | " | 10.0 | | 95.6 | 82-128 | | | | |
| Chloromethane | 9.73 | | " | 10.0 | | 97.3 | 43-155 | | | | |
| cis-1,2-Dichloroethylene | 10.2 | | " | 10.0 | | 102 | 83-129 | | | | |
| cis-1,3-Dichloropropylene | 10.6 | | " | 10.0 | | 106 | 80-131 | | | | |
| Cyclohexane | 5.29 | | " | 10.0 | | 52.9 | 63-149 | Low Bias | | | |
| Dibromochloromethane | 9.16 | | " | 10.0 | | 91.6 | 80-130 | | | | |
| Dibromomethane | 10.2 | | " | 10.0 | | 102 | 72-134 | | | | |
| Dichlorodifluoromethane | 8.50 | | " | 10.0 | | 85.0 | 44-144 | | | | |
| Ethyl Benzene | 12.4 | | " | 10.0 | | 124 | 80-131 | | | | |
| Hexachlorobutadiene | 14.8 | | " | 10.0 | | 148 | 67-146 | High Bias | | | |
| Isopropylbenzene | 12.1 | | " | 10.0 | | 121 | 76-140 | | | | |
| Methyl acetate | 9.25 | | " | 10.0 | | 92.5 | 51-139 | | | | |
| Methyl tert-butyl ether (MTBE) | 8.00 | | " | 10.0 | | 80.0 | 76-135 | | | | |
| Methylcyclohexane | 12.9 | | " | 10.0 | | 129 | 72-143 | | | | |
| Methylene chloride | 10.1 | | " | 10.0 | | 101 | 55-137 | | | | |



Volatile Organic Compounds by GC/MS - Quality Control Data
York Analytical Laboratories, Inc. - Stratford

| Analyte | Result | Reporting Limit | Units | Spike Level | Source* Result | %REC | %REC Limits | Flag | RPD | RPD Limit | Flag |
|---------|--------|-----------------|-------|-------------|----------------|------|-------------|------|-----|-----------|------|
|---------|--------|-----------------|-------|-------------|----------------|------|-------------|------|-----|-----------|------|

Batch BC31487 - EPA 5030B

LCS (BC31487-BS1)

Prepared & Analyzed: 03/21/2023

| | | | | | | | | | | | |
|-----------------------------|------|--|------|------|--|------|--------|--|--|--|--|
| n-Butylbenzene | 12.8 | | ug/L | 10.0 | | 128 | 79-132 | | | | |
| n-Propylbenzene | 12.6 | | " | 10.0 | | 126 | 78-133 | | | | |
| o-Xylene | 12.1 | | " | 10.0 | | 121 | 78-130 | | | | |
| p- & m- Xylenes | 24.6 | | " | 20.0 | | 123 | 77-133 | | | | |
| p-Isopropyltoluene | 12.0 | | " | 10.0 | | 120 | 81-136 | | | | |
| sec-Butylbenzene | 12.7 | | " | 10.0 | | 127 | 79-137 | | | | |
| Styrene | 11.4 | | " | 10.0 | | 114 | 67-132 | | | | |
| tert-Butyl alcohol (TBA) | 21.1 | | " | 50.0 | | 42.1 | 25-162 | | | | |
| tert-Butylbenzene | 10.3 | | " | 10.0 | | 103 | 77-138 | | | | |
| Tetrachloroethylene | 10.3 | | " | 10.0 | | 103 | 82-131 | | | | |
| Toluene | 11.8 | | " | 10.0 | | 118 | 80-127 | | | | |
| trans-1,2-Dichloroethylene | 10.1 | | " | 10.0 | | 101 | 80-132 | | | | |
| trans-1,3-Dichloropropylene | 9.87 | | " | 10.0 | | 98.7 | 78-131 | | | | |
| Trichloroethylene | 11.2 | | " | 10.0 | | 112 | 82-128 | | | | |
| Trichlorofluoromethane | 11.6 | | " | 10.0 | | 116 | 67-139 | | | | |
| Vinyl Chloride | 9.10 | | " | 10.0 | | 91.0 | 58-145 | | | | |

Surrogate: SURR: 1,2-Dichloroethane-d4

9.34

"

10.0

93.4

69-130

Surrogate: SURR: Toluene-d8

10.8

"

10.0

108

81-117

Surrogate: SURR: p-Bromofluorobenzene

9.72

"

10.0

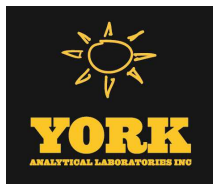
97.2

79-122

LCS Dup (BC31487-BSD1)

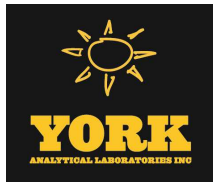
Prepared & Analyzed: 03/21/2023

| | | | | | | | | | | | |
|---|------|--|------|------|--|------|--------|--|-------|----|--|
| 1,1,1,2-Tetrachloroethane | 9.92 | | ug/L | 10.0 | | 99.2 | 82-126 | | 2.00 | 30 | |
| 1,1,1-Trichloroethane | 8.63 | | " | 10.0 | | 86.3 | 78-136 | | 4.64 | 30 | |
| 1,1,2,2-Tetrachloroethane | 11.6 | | " | 10.0 | | 116 | 76-129 | | 4.15 | 30 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 9.85 | | " | 10.0 | | 98.5 | 54-165 | | 5.91 | 30 | |
| 1,1,2-Trichloroethane | 10.5 | | " | 10.0 | | 105 | 82-123 | | 0.287 | 30 | |
| 1,1-Dichloroethane | 9.32 | | " | 10.0 | | 93.2 | 82-129 | | 7.14 | 30 | |
| 1,1-Dichloroethylene | 9.84 | | " | 10.0 | | 98.4 | 68-138 | | 7.91 | 30 | |
| 1,2,3-Trichlorobenzene | 11.3 | | " | 10.0 | | 113 | 76-136 | | 1.96 | 30 | |
| 1,2,3-Trichloropropane | 10.6 | | " | 10.0 | | 106 | 77-128 | | 4.95 | 30 | |
| 1,2,4-Trichlorobenzene | 10.3 | | " | 10.0 | | 103 | 76-137 | | 0.292 | 30 | |
| 1,2,4-Trimethylbenzene | 10.6 | | " | 10.0 | | 106 | 82-132 | | 7.29 | 30 | |
| 1,2-Dibromo-3-chloropropane | 11.1 | | " | 10.0 | | 111 | 45-147 | | 1.45 | 30 | |
| 1,2-Dibromoethane | 10.0 | | " | 10.0 | | 100 | 83-124 | | 3.35 | 30 | |
| 1,2-Dichlorobenzene | 10.7 | | " | 10.0 | | 107 | 79-123 | | 1.85 | 30 | |
| 1,2-Dichloroethane | 9.23 | | " | 10.0 | | 92.3 | 73-132 | | 1.86 | 30 | |
| 1,2-Dichloropropane | 12.2 | | " | 10.0 | | 122 | 78-126 | | 2.59 | 30 | |
| 1,3,5-Trimethylbenzene | 11.0 | | " | 10.0 | | 110 | 80-131 | | 6.42 | 30 | |
| 1,3-Dichlorobenzene | 9.80 | | " | 10.0 | | 98.0 | 86-122 | | 4.20 | 30 | |
| 1,4-Dichlorobenzene | 9.71 | | " | 10.0 | | 97.1 | 85-124 | | 3.04 | 30 | |
| 1,4-Dioxane | 219 | | " | 210 | | 104 | 10-349 | | 10.4 | 30 | |
| 2-Butanone | 9.36 | | " | 10.0 | | 93.6 | 49-152 | | 2.95 | 30 | |
| 2-Hexanone | 10.1 | | " | 10.0 | | 101 | 51-146 | | 7.07 | 30 | |
| 4-Methyl-2-pentanone | 9.33 | | " | 10.0 | | 93.3 | 57-145 | | 6.07 | 30 | |
| Acetone | 9.47 | | " | 10.0 | | 94.7 | 14-150 | | 1.99 | 30 | |
| Acrolein | 5.75 | | " | 10.0 | | 57.5 | 10-153 | | 4.63 | 30 | |
| Acrylonitrile | 8.96 | | " | 10.0 | | 89.6 | 51-150 | | 5.03 | 30 | |
| Benzene | 10.3 | | " | 10.0 | | 103 | 85-126 | | 3.91 | 30 | |
| Bromochloromethane | 10.0 | | " | 10.0 | | 100 | 77-128 | | 0.993 | 30 | |
| Bromodichloromethane | 10.3 | | " | 10.0 | | 103 | 79-128 | | 2.78 | 30 | |



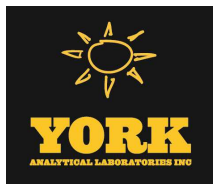
Volatile Organic Compounds by GC/MS - Quality Control Data
York Analytical Laboratories, Inc. - Stratford

| Analyte | Result | Reporting Limit | Units | Spike Level | Source* Result | %REC | %REC Limits | Flag | RPD | RPD Limit | Flag |
|---|--------|-----------------|-------|-------------|----------------|------|-------------|----------|-------|-----------|------|
| Batch BC31487 - EPA 5030B | | | | | | | | | | | |
| LCS Dup (BC31487-BSD1) | | | | | | | | | | | |
| Prepared & Analyzed: 03/21/2023 | | | | | | | | | | | |
| Bromoform | 8.63 | | ug/L | 10.0 | | 86.3 | 78-133 | | 4.62 | 30 | |
| Bromomethane | 7.31 | | " | 10.0 | | 73.1 | 43-168 | | 20.3 | 30 | |
| Carbon disulfide | 9.80 | | " | 10.0 | | 98.0 | 68-146 | | 5.94 | 30 | |
| Carbon tetrachloride | 8.80 | | " | 10.0 | | 88.0 | 77-141 | | 4.77 | 30 | |
| Chlorobenzene | 11.0 | | " | 10.0 | | 110 | 88-120 | | 3.57 | 30 | |
| Chloroethane | 8.29 | | " | 10.0 | | 82.9 | 65-136 | | 8.99 | 30 | |
| Chloroform | 9.24 | | " | 10.0 | | 92.4 | 82-128 | | 3.40 | 30 | |
| Chloromethane | 8.28 | | " | 10.0 | | 82.8 | 43-155 | | 16.1 | 30 | |
| cis-1,2-Dichloroethylene | 9.57 | | " | 10.0 | | 95.7 | 83-129 | | 6.47 | 30 | |
| cis-1,3-Dichloropropylene | 10.4 | | " | 10.0 | | 104 | 80-131 | | 1.33 | 30 | |
| Cyclohexane | 4.95 | | " | 10.0 | | 49.5 | 63-149 | Low Bias | 6.64 | 30 | |
| Dibromochloromethane | 9.41 | | " | 10.0 | | 94.1 | 80-130 | | 2.69 | 30 | |
| Dibromomethane | 10.2 | | " | 10.0 | | 102 | 72-134 | | 0.392 | 30 | |
| Dichlorodifluoromethane | 7.71 | | " | 10.0 | | 77.1 | 44-144 | | 9.75 | 30 | |
| Ethyl Benzene | 11.7 | | " | 10.0 | | 117 | 80-131 | | 5.23 | 30 | |
| Hexachlorobutadiene | 14.1 | | " | 10.0 | | 141 | 67-146 | | 4.71 | 30 | |
| Isopropylbenzene | 11.4 | | " | 10.0 | | 114 | 76-140 | | 6.30 | 30 | |
| Methyl acetate | 9.63 | | " | 10.0 | | 96.3 | 51-139 | | 4.03 | 30 | |
| Methyl tert-butyl ether (MTBE) | 8.59 | | " | 10.0 | | 85.9 | 76-135 | | 7.11 | 30 | |
| Methylcyclohexane | 11.9 | | " | 10.0 | | 119 | 72-143 | | 8.07 | 30 | |
| Methylene chloride | 9.88 | | " | 10.0 | | 98.8 | 55-137 | | 2.10 | 30 | |
| n-Butylbenzene | 11.8 | | " | 10.0 | | 118 | 79-132 | | 7.80 | 30 | |
| n-Propylbenzene | 11.8 | | " | 10.0 | | 118 | 78-133 | | 7.30 | 30 | |
| o-Xylene | 11.6 | | " | 10.0 | | 116 | 78-130 | | 4.47 | 30 | |
| p- & m- Xylenes | 23.3 | | " | 20.0 | | 116 | 77-133 | | 5.40 | 30 | |
| p-Isopropyltoluene | 11.0 | | " | 10.0 | | 110 | 81-136 | | 8.49 | 30 | |
| sec-Butylbenzene | 11.7 | | " | 10.0 | | 117 | 79-137 | | 7.78 | 30 | |
| Styrene | 11.1 | | " | 10.0 | | 111 | 67-132 | | 2.66 | 30 | |
| tert-Butyl alcohol (TBA) | 22.6 | | " | 50.0 | | 45.3 | 25-162 | | 7.23 | 30 | |
| tert-Butylbenzene | 9.61 | | " | 10.0 | | 96.1 | 77-138 | | 7.12 | 30 | |
| Tetrachloroethylene | 9.56 | | " | 10.0 | | 95.6 | 82-131 | | 7.06 | 30 | |
| Toluene | 11.0 | | " | 10.0 | | 110 | 80-127 | | 6.66 | 30 | |
| trans-1,2-Dichloroethylene | 9.76 | | " | 10.0 | | 97.6 | 80-132 | | 3.62 | 30 | |
| trans-1,3-Dichloropropylene | 9.94 | | " | 10.0 | | 99.4 | 78-131 | | 0.707 | 30 | |
| Trichloroethylene | 10.5 | | " | 10.0 | | 105 | 82-128 | | 6.44 | 30 | |
| Trichlorofluoromethane | 10.8 | | " | 10.0 | | 108 | 67-139 | | 6.43 | 30 | |
| Vinyl Chloride | 8.13 | | " | 10.0 | | 81.3 | 58-145 | | 11.3 | 30 | |
| Surrogate: SURRE: 1,2-Dichloroethane-d4 | 9.70 | | " | 10.0 | | 97.0 | 69-130 | | | | |
| Surrogate: SURRE: Toluene-d8 | 10.7 | | " | 10.0 | | 107 | 81-117 | | | | |
| Surrogate: SURRE: p-Bromofluorobenzene | 9.63 | | " | 10.0 | | 96.3 | 79-122 | | | | |



Volatile Analysis Sample Containers

| Lab ID | Client Sample ID | Volatile Sample Container |
|------------|------------------|---|
| 23C0637-01 | DW-1 | 40mL Clear Vial (pre-pres.) HCl; Cool to 4° C |
| 23C0637-02 | DW-2 | 40mL Clear Vial (pre-pres.) HCl; Cool to 4° C |



Sample and Data Qualifiers Relating to This Work Order

| | |
|-------|--|
| QL-02 | This LCS analyte is outside Laboratory Recovery limits due the analyte behavior using the referenced method. The reference method has certain limitations with respect to analytes of this nature. |
| J | Detected below the Reporting Limit but greater than or equal to the Method Detection Limit (MDL/LOD) or in the case of a TIC, the result is an estimated concentration. |
| ICVE | The value reported is ESTIMATED. The value is estimated due to its behavior during initial calibration verification (recovery exceeded 30% of expected value). |
| CCVE | The value reported is ESTIMATED. The value is estimated due to its behavior during continuing calibration verification (>20% Difference for average Rf or >20% Drift for quadratic fit). |

Definitions and Other Explanations

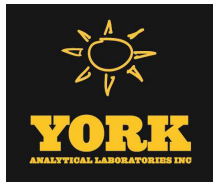
| | |
|-------------|--|
| * | Analyte is not certified or the state of the samples origination does not offer certification for the Analyte. |
| ND | NOT DETECTED - the analyte is not detected at the Reported to level (LOQ/RL or LOD/MDL) |
| RL | REPORTING LIMIT - the minimum reportable value based upon the lowest point in the analyte calibration curve. |
| LOQ | LIMIT OF QUANTITATION - the minimum concentration of a target analyte that can be reported within a specified degree of confidence. This is the lowest point in an analyte calibration curve that has been subjected to all steps of the processing/analysis and verified to meet defined criteria. This is based upon NELAC 2009 Standards and applies to all analyses. |
| LOD | LIMIT OF DETECTION - a verified estimate of the minimum concentration of a substance in a given matrix that an analytical process can reliably detect. This is based upon NELAC 2009 Standards and applies to all analyses conducted under the auspices of EPA SW-846. |
| MDL | METHOD DETECTION LIMIT - a statistically derived estimate of the minimum amount of a substance an analytical system can reliably detect with a 99% confidence that the concentration of the substance is greater than zero. This is based upon 40 CFR Part 136 Appendix B and applies only to EPA 600 and 200 series methods. |
| Reported to | This indicates that the data for a particular analysis is reported to either the LOD/MDL, or the LOQ/RL. In cases where the "Reported to" is located above the LOD/MDL, any value between this and the LOQ represents an estimated value which is "J" flagged accordingly. This applies to volatile and semi-volatile target compounds only. |
| NR | Not reported |
| RPD | Relative Percent Difference |
| Wet | The data has been reported on an as-received (wet weight) basis |
| Low Bias | Low Bias flag indicates that the recovery of the flagged analyte is below the laboratory or regulatory lower control limit. The data user should take note that this analyte may be biased low but should evaluate multiple lines of evidence including the LCS and site-specific MS/MSD data to draw bias conclusions. In cases where no site-specific MS/MSD was requested, only the LCS data can be used to evaluate such bias. |
| High Bias | High Bias flag indicates that the recovery of the flagged analyte is above the laboratory or regulatory upper control limit. The data user should take note that this analyte may be biased high but should evaluate multiple lines of evidence including the LCS and site-specific MS/MSD data to draw bias conclusions. In cases where no site-specific MS/MSD was requested, only the LCS data can be used to evaluate such bias. |
| Non-Dir. | Non-dir. flag (Non-Directional Bias) indicates that the Relative Percent Difference (RPD) (a measure of precision) among the MS and MSD data is outside the laboratory or regulatory control limit. This alerts the data user where the MS and MSD are from site-specific samples that the RPD is high due to either non-homogeneous distribution of target analyte between the MS/MSD or indicates poor reproducibility for other reasons. |

If EPA SW-846 method 8270 is included herein it is noted that the target compound N-nitrosodiphenylamine (NDPA) decomposes in the gas chromatographic inlet and cannot be separated from diphenylamine (DPA). These results could actually represent 100% DPA, 100% NDPA or some combination of the two. For this reason, York reports the combined result for n-nitrosodiphenylamine and diphenylamine for either of these compounds as a combined concentration as Diphenylamine.

If Total PCBs are detected and the target aroclors reported are "Not detected", the Total PCB value is reported due to the presence of either or both Aroclors 1262 and 1268 which are non-target aroclors for some regulatory lists.

2-chloroethylvinyl ether readily breaks down under acidic conditions. Samples that are acid preserved, including standards will exhibit breakdown. The data user should take note.

Certification for pH is no longer offered by NYDOH ELAP.



Semi-Volatile and Volatile analyses are reported down to the LOD/MDL, with values between the LOD/MDL and the LOQ being "J" flagged as estimated results.

For analyses by EPA SW-846-8270D, the Limit of Quantitation (LOQ) reported for benzidine is based upon the lowest standard used for calibration and is not a verified LOQ due to this compound's propensity for oxidative losses during extraction/concentration procedures and non-reproducible chromatographic performance.




York Analytical Laboratories, Inc.
 120 Research Drive
 Stratford, CT 06615
 clientservices@yorklab.com
 www.yorklab.com

Field Chain-of-Custody Record

YORK Project No.
 2300037

Page of

NOTE: YORK's Standard Terms & Conditions are listed on the back side of this document. This document serves as your written authorization for YORK to proceed with the analyses requested below. Your signature binds you to YORK's Standard Terms & Conditions.

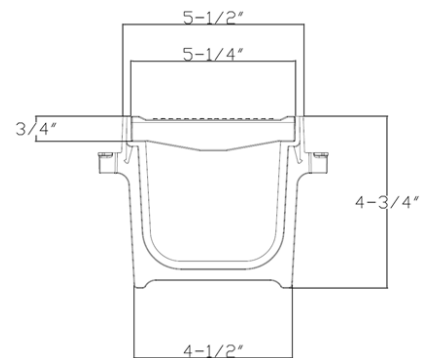
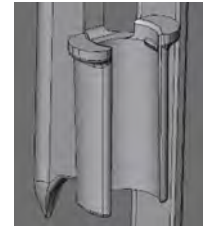
| YOUR INFORMATION | | Report To: | | Invoice To: | | YOUR Project Number | | Turn-Around Time | | |
|---|-------------------------------|---|--|---|--|--|--------------------|-----------------------|-------------------------------------|---------------------------|
| Company: | Advanced Cleanup Technologies | Company: | ACT | Company: | ACT | 9628 - BXNY | | RUSH - Next Day | <input type="checkbox"/> | |
| Address: | 235 Park Ave PMB 34864 | Address: | Same | Address: | Same | YOUR Project Name | | RUSH - Two Day | <input type="checkbox"/> | |
| Phone: | NY, NY 10003 | Phone: | | Phone: | | | | RUSH - Three Day | <input type="checkbox"/> | |
| Contact: | Jasen Stewart | Contact: | Paul Stewart | Contact: | Karen Friedman | | | RUSH - Four Day | <input type="checkbox"/> | |
| E-mail: | jasen@act.earth | E-mail: | PaulStewart@act.earth | E-mail: | KarenF@act.earth | YOUR PO#: | | Standard (5-7 Day) | <input checked="" type="checkbox"/> | |
| Please print clearly and legibly. All information must be complete. Samples will not be logged in and the turn-around-time clock will not begin until any questions by YORK are resolved. | | | | | | | | | | |
| Samples Collected by: (print your name above and sign below)  | | | | | | | | | | |
| Sample Identification | | Matrix Codes | Report / EDD Type (circle selections) | YORK Reg. Comp. | Samples From | | Analysis Requested | Container Description | Special Instruction | |
| DW-1 | DW-2 | S - soil / solid GW - groundwater DW - drinking water WW - wastewater O - Oil ; Other | CT RCP CT RCP DQA/DUE NJDEP Reduced Deliverables NJDKQP | Compared to the following Regulation(s): (please fill in) NYSDEC part 375 TGS(GA) | New York New Jersey Connecticut Pennsylvania Other | Standard Excel EDD EQUS (Standard) NJDEP SRP HazSite Other: | VOL " | 3 VOLs HCL " | Field Filtered Lab to Filter | Date/Time 3/10/23 1700 |
| Date/Time Sampled | | Sample Matrix | Preservation: (check all that apply) | Samples Received by / Company | | Samples Relinquished by / Company | | Temp. Received at Lab | | Degrees C |
| 3/9/23 1320 " 1245 | | OW " | HCl <input checked="" type="checkbox"/> MeOH <input type="checkbox"/> HNO3 <input type="checkbox"/> H2SO4 <input type="checkbox"/> NaOH <input type="checkbox"/> ZnAc <input type="checkbox"/> Ascorbic Acid <input type="checkbox"/> Other: <input type="checkbox"/> | KBarhyork 920 AM 3/10/23 | | KBarhyork 920 AM 3/10/23 | | 1700 | | 3.0 |
| Comments: | | | | | | | | | | |
| ACT BRIDGE | | | | | | | | | | |

**Appendix F:
Engineering Controls
Datasheets**

5 Inch Pro Series Drain Kit



NDS #864
Polypropylene



Part #: 864 (Includes #814-Grates (2qty.), #800-Channel (1qty.), #813-End Outlet (1qty.), and #812-End Cap (1qty.)

Material: Channel (Polyolefin) Grates (HDPE)

Color: Light Gray

Fits: 3" (Hub) and 4" (Spigot) Sewer/Drain Pipe

Rebar tie clips for easier installation: Fits #4 Rebar

Grate Opening: 3/8" x 1-1/4"

Open Surface Area: 23.52 Sq. Inch per Ft.

Head Pressure / Flow Rate:

Head (inches) - Max Flow

1" = 101.75 GPM

0.5" = 71.95 GPM

Weight per each: 5.67 lbs.

Screw: #829 Stainless Steel Screw, 4 per grate.

UV Inhibitors



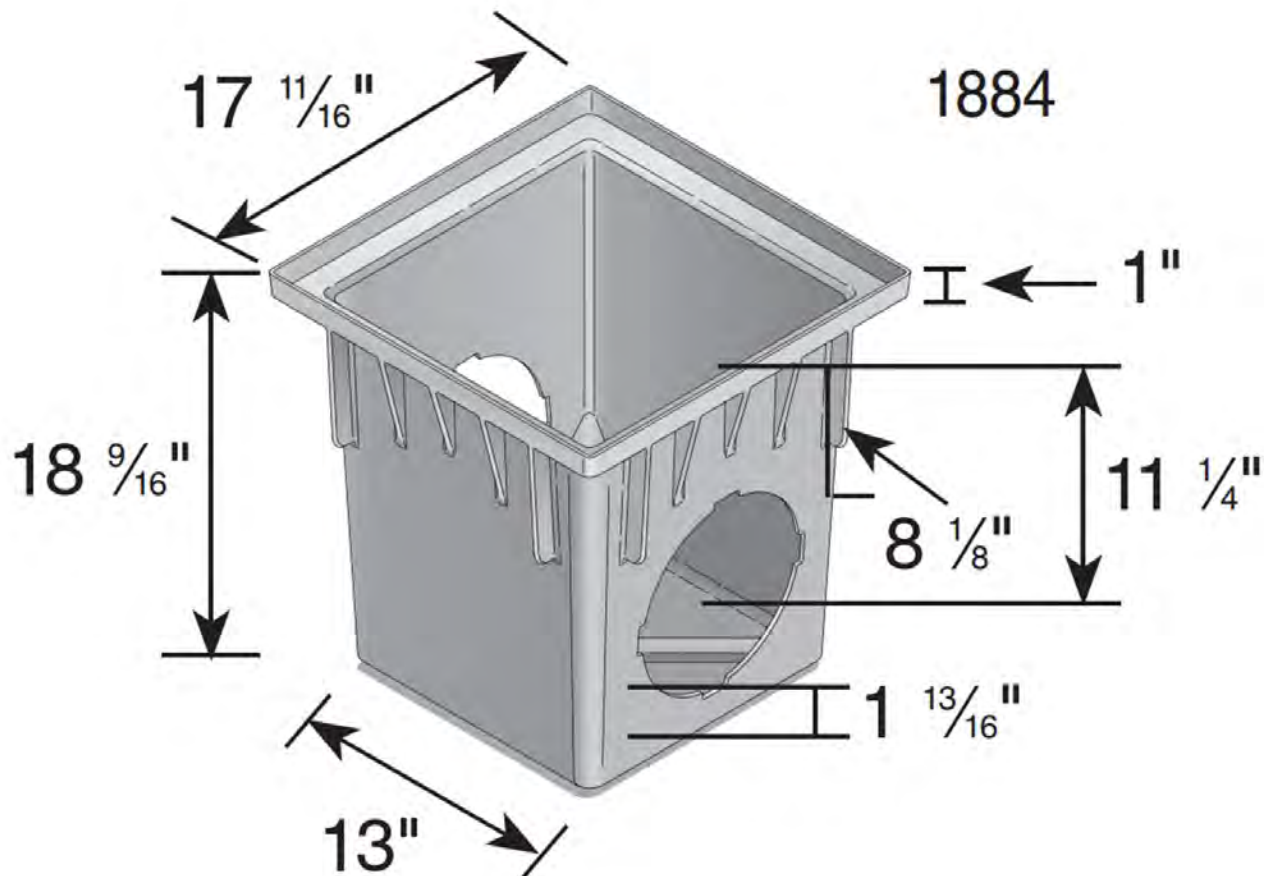
Class B

- Loads of 61-175 psi.
- Recommended for medium-duty pneumatic tire traffic, autos and light trucks at speeds less than 20 m.p.h.

ADA Compliant



18" ONE PIECE CATCH BASIN



Material: UV Protected High Density Polyethylene (HDPE)

Weight: 9.65 lbs

Colors: Black

- 18" One Piece Catch Basin with 2 Openings (1882)
Includes 2 #1890 Reducer Rings
- 18" One Piece Catch Basin with 4 Openings (1884)
Includes 4 #1890 Reducer Rings

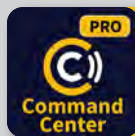
Requires either #1206, #1242, #1245, #1266, #1888, or #1889 universal outlet to connect pipe



POWERFUL AND QUIET AIR FILTRATION HEPA 700

PORTABLE HIGH PERFORMANCE
AIR PURIFICATION FOR WATER AND
FIRE DAMAGE RESTORATION, MOLD
REMEDICATION AND CONSTRUCTION

- **Quieter operation** for use in any occupied spaces
- **Filters out 99.99% of dust and mold** with prefilter and true HEPA filter media - passes DOP tests with no additional filters
- **Removes smoke, odors and fumes fast** with optional carbon filter
- **Lightweight and stacks securely** for storage and transport
- **Variable speed** enables fine-tuned airflow for containments or ducted operation
- **24/7 remote monitoring and control** using Command Center app to ensure continuous air filtration



**NEW COMMAND HUB
FOR REMOTE OPERATION
FROM ANYWHERE!**



FEATURES AND SPECS:

- Variable to 700 CFM with up to 6 ACH in 7000 cu. ft. spaces
- Digital touchscreen displays CFM for better airflow control
- Low amp draw - 1.5 - 3 amps
- Convenient onboard GFCI outlet for connecting up to 8 units on 15-amp circuit
- Rugged injection-molded housing with limited lifetime warranty



| | |
|--------------------|--|
| MODEL | HEPA 700 (#125105) |
| AIRFLOW | VARIABLE UP TO 700 CFM |
| STATIC PRESSURE | 3.0 IN. 7.6 CM (H2O) |
| WEIGHT | 42 LBS. 19 KG |
| DIMENSIONS (HxDxW) | 23 X 17 X 26 IN. 57.9 X 43.3 X 65.9 CM |
| POWER (115V, 60HZ) | 1.5 - 3 AMPS |
| DUCT SIZE | 14 IN. 35.6 CM |
| XACTIMATE | WTRNAFAN> |

Final specifications are subject to change.



**Appendix G:
Data Usability
Summary Reports**

December 8, 2025

Mr. Jason Stewart
Advanced Cleanup Technologies
228 Park Ave S PMB 34864
New York, New York 10003

Re: Data Usability Summary Report – York Analytical Laboratories, Inc. - 20H0017

Dear Mr. Stewart:

The evaluation of analytical data by York Analytical Laboratories for project 9268-BXNY, which were reported in a single data package under Job No. 20H0017 has been completed. The following samples were reported.

| | | | |
|------|------|------|------|
| SS-1 | SS-2 | IA-1 | SS-3 |
| SS-4 | IA-2 | OA-1 | |

Analysis was performed in accordance with EPA Method TO-15 (volatile organics). The review was performed to the extent possible, in accordance with the analytical method, and “DER-10/ Technical Guidance for Site Investigation and Remediation”. Professional judgment is applied as necessary and appropriate. National Functional Guidelines for Organic Data Review was consulted as needed. Qualifiers consistent with those defined by EPA Region 2 are applied as necessary and appropriate.

Below is the Data Usability Summary Report (DUSR) associated with these samples.

| Data Usability Summary Report | |
|---|---|
| 1. Is the data package complete as defined under the requirements for the most current DEC ASP Category B or USEPA CLP data deliverables? | No -see Documentation section and section A.3 |
| 2. Have all holding times been met? | Yes |
| 3. Do all the QC data; blanks, instrument tunings, calibration standards, calibration verifications, surrogate recoveries, spike recoveries, replicate analyses, laboratory controls and sample data fall within the protocol required limits and specifications? | No -see following sections |
| 4. Have all of the data been generated using established and agreed upon analytical protocols? | Yes |
| 5. Does an evaluation of the raw data confirm the results provided in the data summary sheet and the quality control verification forms? | Yes |
| 6. Have the correct data qualifiers been used and are they consistent with the most current DEC ASP? | Yes |

| | |
|---|-----|
| 7. Have any quality control (QC) exceedances been specifically noted in the DUSR and have the corresponding QC summary sheet from the data package been attached to the DUSR? | Yes |
|---|-----|

Overall Evaluation

Based on the data review effort, results are usable, with the following qualifications. For samples that are qualified as estimated (J-, UJ), detected results may be biased low. False negatives may exist in non-detect results. Sample results that are qualified as estimated (J+) may be biased high. For samples that are qualified as estimated with any combination of (J), (J-) and/or (J+), the (J) qualifier takes precedence and is applied to the sample result. It is not possible to determine the direction of the bias and the overall effect on the result.

- The results for all target analytes in all samples are qualified as estimated (J, UJ) because there is no documentation provided to show samples were properly preserved.
- The results for all target analytes in SS-1 are qualified as estimated (J, UJ) due to possible leak during sampling.
- The results for benzyl chloride in all samples are qualified as estimated (UJ) due to high relative standard deviation (RSD) in the initial calibration (IC).
- Based on professional judgment, the results for acetone in SS-1, SS-2, SS-3, SS-4, IA-2 and OA-1 and isopropanol in IA-2 and OA-1 are qualified as estimated (J) due to unacceptable initial calibration.
- The results for 1,4-dioxane and 4-methyl-2-pentanone in all samples are qualified as estimated (UJ) due to low response in the continuing calibration verification (CCV) standard.
- The results for 1,4-dioxane and 4-methyl-2-pentanone in all samples are qualified as estimated (J-, UJ) due to low recovery in the laboratory control sample (LCS).
- Based on professional judgment, samples results less than ten times the quantitation limit in all samples are qualified as estimated (J) because proper canister certification cannot be confirmed.
- The result for isopropanol in IA-1 is qualified as estimated (J) because the reported concentration is outside the calibration range.

Qualifier definitions are provided in Attachment A. A copy of the chain of custody record is provided in Attachment B. Pages from the data package illustrating the exceedances and issues described in this validation report are included in Attachment C. Annotated Summary Forms are included in Attachment D detailing qualifications resulting from the data review effort.

The following components were reviewed, where applicable:

- Chain of Custody
- Receiving conditions

- Holding times
- Preservation
- Analyte lists
- Reporting limits
- Requested methods
- Units, and
- Sample related quality control data:
 - Method, instrument blanks
 - Clean canister certification
 - Field blanks
 - Surrogate recoveries
 - LCS recoveries
 - Internal standard area response
 - Duplicates
 - Analyte Identification
- Instrument related quality control data:
 - Instrument tunes
 - Calibration summaries

The following sections of the report detail only quality control exceedances that impacted results. Where a quality control item exceeded control limits but there is no impact to the samples results, these are not detailed in the report.

Documentation: A completeness review of the data package was performed, and the data package was determined to be a complete Category B data package. The following documentation issues were observed during the review:

- The data package indicates that the laboratory did not hold certification for ethyl acetate, p-ethyltoluene, propylene, tetrahydrofuran, 1,1,1,2-tetrachloroethane, 1,3-dichloropropane, and 2-hexanone. It is noted that NY ELAP does not offer certification for ethyl acetate, p-ethyltoluene, propylene, tetrahydrofuran and 2-hexanone, but does offer certification for 1,1,1,2-tetrachloroethane and 1,3-dichloropropane. The laboratory was contacted to request clarification. The laboratory stated *'Given the age of the project, it is difficult to ascertain the exact circumstances in which these compounds were requested. We don't typically report non-certified compounds, except at the request of the client, so it can be safely assumed that there was correspondence regarding the accreditation status of the compounds at the time the samples were taken. Just to note- while these compounds are indeed non-certified under our scope of accreditation between 2020-2022, we ran the compounds in the exact same way we ran our certified compounds – quantitated against a multiple calibration, where they were properly qualified as specified by our certifying body.'*
- The data package does not include the clean canister certification data. The laboratory was contacted and was requested to provide the missing documentation. The laboratory stated *'our standard archival period is three years and due to this data*

being older than that we won't be able to obtain the additional documents.' See Section A.3.

Holding Times, Preservation, Sample Integrity:

A copy of the applicable chain of custody (COC) record was included in the data package, documenting sample collection date of July 31, 2020. The samples were received at the laboratory on the same day on August 3, 2020. The samples were analyzed within method hold time. However, the data package does contain sample receipt information regarding condition of canisters upon receipt. The laboratory was contacted for clarification and stated *'Any anomalous receipt data would be denoted in the case narrative. If there was nothing listed there wasn't any anomalous activity, but we've opened a corrective action to improve this process.'* Since there is no evidence that the samples were properly preserved (pressure inside the canister maintained within +/- 5 psi from sampling to check in the laboratory or analysis), the results for all target analytes in all samples are qualified as estimated (J, UJ).

The COC indicates that the 'canister gauge didn't move much' for sample SS-1. The initial canister vacuum is recorded at -30mm Hg and the final canister vacuum at -22 mmHg, indicating very little sample was collected and the possibility of a leak existed. On this basis, the results for all target analytes in SS-1 are qualified as estimated (J, UJ).

A. Volatile Organics

1. Calibration

One initial calibration (IC) was performed in support of the sample analysis. All relative response factors (RRFs) and relative standard deviations (RSDs) or correlation coefficients are acceptable with the exceptions noted below.

| Analyte | RSD |
|-----------------|------------|
| Acetone | 100.5 |
| Benzyl chloride | 32.3 |
| Isopropanol | 82.1 |

The laboratory applied quadratic equations to achieve an acceptable correlation coefficient and used this model to quantitate sample concentrations. However, a review of the individual calibration levels indicates a severe lack of instrument sensitivity in the lower calibration standards as presented below.

| Analyte | RRF 50 | RRF 30 | RRF 20 | RRF 10 | RRF 3 | RRF 0.5 | RRF 0.2 | RRF 0.1 |
|----------------|---------------|---------------|---------------|---------------|--------------|----------------|-----------------|-----------------|
| Acetone | 1.171301 | 1.16144 | 1.19475 | 1.22118 | 1.386964 | 4.44930 | 6.520248 | 10.42493 |

| | | | | | | | | |
|-----------------|----------|----------|---------|---------|----------|----------------|-----------------|-----------------|
| Benzyl chloride | 0.211000 | 0.207986 | 0.20257 | 0.20066 | 0.197759 | 0.11883 | 0.096370 | 0.09877 |
| Isopropanol | 1.332178 | 1.318993 | 1.36216 | 1.40152 | 1.693545 | 4.94095 | 6.122283 | 8.010574 |

The high RSD (32.3) for benzyl chloride is caused by the low responses in the two lowest calibration standards. Benzyl chloride is not detected in any samples. The results for benzyl chloride in all samples are qualified as estimated (UJ) due to high RSD in the initial calibration.

The high RSDs for acetone and isopropanol are caused by excessively high response factors in the three lowest concentration standards. The validator calculated the RSD with these three standards excluded, and the RSDs are:

- Acetone: 7.52% RSD, average relative response factor = 1.22711.
- Isopropanol: 39.1% RSD, average relative response factor=1.421679.

The recalculated sample concentrations using the updated average relative response factors are presented below:

| Sample | Reported Acetone (on-column ppbv) | Recalculated Acetone (on-column ppbv) | %D |
|--------|---------------------------------------|---|------|
| SS-1 | 5.79 | 6.18 | 6.7 |
| SS-2 | 3.91 | 4.47 | 14.3 |
| IA-1 | 42.22 | 40.21 | 4.8 |
| SS-3 | 5.18 | 5.62 | 8.5 |
| SS-4 | 2.18 | 2.91 | 33.5 |
| IA-2 | 4.46 | 4.97 | 11.4 |
| OA-1 | 3.97 | 4.5 | 13.4 |
| Sample | Reported Isopropanol (on-column ppbv) | Recalculated Isopropanol (on-column ppbv) | %D |
| SS-1 | ND | NA | NA |
| SS-2 | ND | NA | NA |
| IA-1 | 213.18 (E) | 214.2 | 0.5 |
| SS-3 | ND | NA | NA |
| SS-4 | ND | NA | NA |
| IA-2 | 1.19 | 1.91 | 60.5 |
| OA-1 | 2.80 | 3.36 | 20.0 |

ND-not detected

NA-not applicable

Where the difference in concentration is less than 5%, no action is taken. Where the difference in concentration exceeds 5% sample results are qualified as estimated (J-), biased low due to unacceptable calibration, as follows:

- Acetone in SS-1, SS-2, SS-3, SS-4, IA-2, and OA-1.
- Isopropanol in IA-2 and OA-1.

A second source ICV standard was analyzed after the IC, and all percent differences are acceptable ($\leq 30\%D$). CCVs were analyzed at the appropriate frequency and are acceptable ($\%D \leq 30$) with the exceptions noted below.

| Analyte | %D | Associated Sample | Qualifier Applied |
|---------------------------|-------|--|-------------------|
| <i>CCV 07/07/20 18:12</i> | | | |
| 1,4-Dioxane | -32.9 | SS-1 | J, UJ |
| 4-Methyl-2-pentanone | -31.1 | SS-2 IA-1 SS-3 SS-4 IA-2 OA-1 | |

The percent differences represent a decrease in instrument sensitivity. The results for 1,4-dioxane and 4-methyl-2-pentanone in all samples are qualified as estimated (UJ) due to low response in the CCV.

2. Laboratory Control Sample (LCS) /LCS Duplicate (LCSD)

One LCS was prepared and analyzed with the sample. All recoveries are acceptable (70-130%R), with the following exceptions:

| Analyte | LCS %R | Affected Sample | Qualifier Applied |
|----------------------|--------|--|-------------------|
| <i>BH00312-BS1</i> | | | |
| 1,4-Dioxane | 65.3 | SS-1 | J-, UJ |
| 4-Methyl-2-pentanone | 68.0 | SS-2 IA-1 SS-3 SS-4 IA-2 OA-1 | |

The results for 1,4-dioxane and 4-methyl-2-pentanone in all samples are qualified as estimated (J-, UJ) due to low recovery in the LCS.

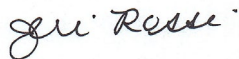
3. Compound Quantitation

As noted in **Documentation** section, the laboratory did not provide clean canister certification data. Since clean canister certification data is not available, it cannot be confirmed whether the canisters used for sampling were properly cleaned and free of background contamination. This uncertainty introduces potential impacts to sample results, including the possibility that detected compounds may originate from residual contamination within the canisters rather than the sampled environment. As a result, accuracy cannot be fully assured, and reported concentrations may not reliably represent actual ambient conditions. Based on professional judgment, samples results less than ten times the quantitation limit in all samples are qualified as estimated (J) because proper canister certification cannot be confirmed.

The result for isopropanol in IA-1 is qualified as estimated (J) because the reported concentration is outside the calibration range.

No other sample results are qualified. Please feel free to contact me at (908) 370-3431 or richjerirossi513@gmail.com if you have any questions regarding this data package review report or need further information.

Sincerely,

A handwritten signature in cursive script that reads "Jeri Rossi".

Jeri L Rossi, CEAC

Environmental Consulting Chemist

ATTACHMENT A

Qualifier Definitions

EPA Qualifier Definitions

- U The analyte was analyzed for but was not detected above the level of the reported sample quantitation limit.
- J The result is an estimated quantity. The associated numerical value is the approximate concentration of the analyte in the sample.
- J+ The result is an estimated quantity, but the result may be biased high.
- J- The result is an estimated quantity, but the result may be biased low.
- NJ The analyte has been “tentatively identified” or “presumptively” as present and the associated numerical value is the estimated concentration in the sample.
- UJ The analyte was analyzed for but was not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
- R The data are unusable. The sample results are rejected due to serious deficiencies in meeting QC criteria. The analyte may or may not be present in the sample.

ATTACHMENT B

CHAIN OF CUSTODY (COC)



York Analytical Laboratories, Inc.
120 Research Drive Stratford, CT 06615
132-02 89th Ave Queens, NY 11418

Field Chain-of-Custody Record - AIR

YORK Project No.

20110017



clientservices@yorklab.com
www.yorklab.com

NOTE: YORK's Standard Terms & Conditions are listed on the back side of this document. This document serves as your written authorization for YORK to proceed with the analyses requested below. signature binds you to YORK's Standard Terms & Conditions.

Your Page ___ of ___

| | | | | | | | | | |
|--|--|--------------------------------|--|----------------------------------|--|----------------------------|--|--|--|
| YOUR Information | | Report To: | | Invoice To: | | YOUR Project Number | | Turn-Around Time | |
| Company: <u>Advanced Cleanup Tech</u> | | Company: <u>ACT</u> | | Company: <u>ACT</u> | | <u>9628-BXNY</u> | | RUSH - Next Day | |
| Address: <u>200 Broadhollow Rd. Melville, NY</u> | | Address: <u>Same</u> | | Address: <u>Same</u> | | | | RUSH - Two Day | |
| Phone: <u>516-441-5300</u> | | Phone: _____ | | Phone: _____ | | YOUR Project Name | | RUSH - Three Day | |
| Contact: <u>Chelsea Ferrucci</u> | | Contact: <u>Paul Stewart</u> | | Contact: <u>Karen Friedman</u> | | | | RUSH - Four Day | |
| E-mail: <u>Chelsea.F@act-earth</u> | | E-mail: <u>pauls@act-earth</u> | | E-mail: <u>Karen.F@act-earth</u> | | YOUR PO#: | | Standard (5-7 Day) <input checked="" type="checkbox"/> | |

| | | | | | | |
|---|--|---------------------|--|------------------|----------------------|---|
| Please print clearly and legibly. All information must be complete. Samples will not be logged in and the turn-around-time clock will not begin until any questions by YORK are resolved. | Air Matrix Codes | Samples From | Report / EDD Type (circle selections) | | | YORK Reg. Comp. Compared to the following Regulation(s): (please fill in) |
| | AI - Indoor Ambient Air | New York | <input checked="" type="checkbox"/> | Summary Report | CT RCP | |
| Samples Collected by: (print your name above and sign below) <u>Tim Yang</u> <u>T-Y</u> | AO - Outdoor Amb. Air | New Jersey | <input type="checkbox"/> | QA Report | CT RCP DQA/DUE | EQUIS (Standard) |
| | AE - Vapor Extraction Well/ Process Gas/Effluent | Connecticut | <input type="checkbox"/> | NY ASP A Package | NJDEP Reduced Deliv. | NYSDEC EQUIS |
| | AS - Soil Vapor/Sub-Slab | Pennsylvania | <input type="checkbox"/> | NY ASP B Package | NJDQKP | NJDEP SRP HazSite |
| | | Other | <input type="checkbox"/> | Other: | | |

| Certified Canisters: Batch ___ Individual ___ | | Please enter the following REQUIRED Field Data | | | | | | Reporting Units: ug/m ³ ___ ppbv ___ ppmv ___ | |
|---|-------------------|--|---|--|-------------|---------------|--------------------|--|--|
| Sample Identification | Date/Time Sampled | Air Matrix | Canister Vacuum Before Sampling (in Hg) | Canister Vacuum After Sampling (in Hg) | Canister ID | Flow Cont. ID | Analysis Requested | | |
| SS-1 | 7/31/20 10-1400 | AS | -30 | -22 | 28303 | 7087 | * TO-15 | | |
| SS-2 | " " | " | -30 | -12 | 23197 | 6873 | " | | |
| IA-1 | " " | AI | -30 | -8 | 34496 | 5116 | " | | |
| SS-3 | " " | AS | -30 | -8 | 24128 | 7607 | " | | |
| SS-4 | " " | " | -30 | -8 | Y61 | Y30 | " | | |
| IA-2 | " " | AI | -30 | -8 | 23993 | 7417 | " | | |
| OA-1 | " " | AO | -30 | -8 | 23800 | 5705 | " | | |

| | | | |
|---|----------------------------------|----------------------|--|
| Comments: <u>* canister gauge didnt move much - ran for 5 hrs -</u> | Detection Limits Required | | Sampling Media |
| | ≤ 1 ug/m ³ ___ | NYSDEC V1 Limits ___ | 6 Liter Canister <input checked="" type="checkbox"/> |
| | Routine Survey ___ | Other ___ | Tedlar Bag |

| | | | | | |
|-----------------------------------|-----------------------|-----------------------------------|-----------------------|-----------------------------------|------------------------|
| Samples Relinquished by / Company | Date/Time | Samples Received by / Company | Date/Time | Samples Relinquished by / Company | Date/Time |
| <u>T-Y ACT</u> | <u>8/3/20 9:50 AM</u> | <u>KBaduyok</u> | <u>8/3/20 9:50 AM</u> | <u>KBaduyok</u> | <u>8/3/20 1615</u> |
| Samples Received by / Company | Date/Time | Samples Relinquished by / Company | Date/Time | Samples Received by / Company | Date/Time |
| <u>TC Yalib / YORK</u> | <u>8/3/20 1615</u> | <u>TC Yalib / YORK</u> | <u>8/3/20 1955</u> | <u>Ed / York</u> | <u>8/3/20 1955</u> |
| Samples Relinquished by / Company | Date/Time | Samples Received by / Company | Date/Time | Samples Received in LAB by | Date/Time |
| <u>Ed / York</u> | <u>8/3/20 2150</u> | <u>Lab</u> | <u>8/3/20 2200</u> | <u>A</u> | <u>8/6/20 11:40 PM</u> |

Page 40 of 311

ATTACHMENT C

**SELECTED PAGES FROM DATA PACKAGE –
QC EXCEEDANCES AND VALIDATION ISSUES**

FORM III

LCS / LCS DUPLICATE RECOVERY

EPA TO-15

Laboratory: York Analytical Laboratories, Inc. - StratfordSDG: 20H0017Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYMatrix: AirBatch: BH00312Laboratory ID: BH00312-BS1Preparation: EPA TO15 PREPInitial/Final: 400 mL / 400 mL

| COMPOUND | SPIKE ADDED (ppbv) | LCS CONCENTRATION (ppbv) | LCS % REC. # | QC LIMITS REC. |
|---|--------------------|--------------------------|--------------|----------------|
| 1,1,1,2-Tetrachloroethane | 10.0 | 9.40 | 94.0 | 70 - 130 |
| 1,1,1-Trichloroethane | 10.0 | 9.82 | 98.2 | 70 - 130 |
| 1,1,2,2-Tetrachloroethane | 10.0 | 8.98 | 89.8 | 70 - 130 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 10.0 | 10.2 | 102 | 70 - 130 |
| 1,1,2-Trichloroethane | 10.0 | 8.44 | 84.4 | 70 - 130 |
| 1,1-Dichloroethane | 10.0 | 9.83 | 98.3 | 70 - 130 |
| 1,1-Dichloroethylene | 10.0 | 9.42 | 94.2 | 70 - 130 |
| 1,2,4-Trichlorobenzene | 10.0 | 9.97 | 99.7 | 70 - 130 |
| 1,2,4-Trimethylbenzene | 10.0 | 8.49 | 84.9 | 70 - 130 |
| 1,2-Dibromoethane | 10.0 | 8.54 | 85.4 | 70 - 130 |
| 1,2-Dichlorobenzene | 10.0 | 10.1 | 101 | 70 - 130 |
| 1,2-Dichloroethane | 10.0 | 8.86 | 88.6 | 70 - 130 |
| 1,2-Dichloropropane | 10.0 | 7.81 | 78.1 | 70 - 130 |
| 1,2-Dichlorotetrafluoroethane | 10.0 | 10.2 | 102 | 70 - 130 |
| 1,3,5-Trimethylbenzene | 10.0 | 8.97 | 89.7 | 70 - 130 |
| 1,3-Butadiene | 10.0 | 10.8 | 108 | 70 - 130 |
| 1,3-Dichlorobenzene | 10.0 | 10.4 | 104 | 70 - 130 |
| 1,3-Dichloropropane | 10.0 | 7.91 | 79.1 | 70 - 130 |
| 1,4-Dichlorobenzene | 10.0 | 10.4 | 104 | 70 - 130 |
| 1,4-Dioxane | 10.0 | 6.53 | 65.3 * | 70 - 130 |
| 2-Butanone | 10.0 | 9.16 | 91.6 | 70 - 130 |
| 2-Hexanone | 10.0 | 7.34 | 73.4 | 70 - 130 |
| 3-Chloropropene | 10.0 | 10.0 | 100 | 70 - 130 |
| 4-Methyl-2-pentanone | 10.0 | 6.80 | 68.0 * | 70 - 130 |
| Acetone | 10.0 | 9.23 | 92.3 | 70 - 130 |
| Acrylonitrile | 10.0 | 9.82 | 98.2 | 70 - 130 |
| Benzene | 10.0 | 9.79 | 97.9 | 70 - 130 |
| Benzyl chloride | 10.0 | 9.97 | 99.7 | 70 - 130 |
| Bromodichloromethane | 10.0 | 7.86 | 78.6 | 70 - 130 |
| Bromoform | 10.0 | 10.6 | 106 | 70 - 130 |

FORM VI

INITIAL CALIBRATION DATA

EPA TO-15

Laboratory: York Analytical Laboratories, Inc. - StratfordSDG: 20H0017Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YG00006Instrument: 5975CCalibration Date: 07/07/20 18:12

| Compound | Level 01 | | Level 02 | | Level 03 | | Level 04 | | Level 05 | | Level 06 | |
|---|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| 1,1,1,2-Tetrachloroethane | 50 | 0.4403751 | 30 | 0.4484318 | 20 | 0.4594578 | 10 | 0.4855418 | 3 | 0.5696907 | 0.5 | 0.4747171 |
| 1,1,1-Trichloroethane | 50 | 2.519017 | 30 | 2.524141 | 20 | 2.618182 | 10 | 2.708008 | 3 | 3.214963 | 0.5 | 2.630602 |
| 1,1,2,2-Tetrachloroethane | 50 | 0.673399 | 30 | 0.6856557 | 20 | 0.6991994 | 10 | 0.7415378 | 3 | 0.8845473 | 0.5 | 0.711751 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 50 | 2.229221 | 30 | 2.225928 | 20 | 2.321697 | 10 | 2.40721 | 3 | 2.870833 | 0.5 | 2.376132 |
| 1,1,2-Trichloroethane | 50 | 0.3073361 | 30 | 0.3074599 | 20 | 0.3150248 | 10 | 0.3261175 | 3 | 0.3887162 | 0.5 | 0.3145029 |
| 1,1-Dichloroethane | 50 | 1.792969 | 30 | 1.785278 | 20 | 1.845791 | 10 | 1.911867 | 3 | 2.266373 | 0.5 | 1.901432 |
| 1,1-Dichloroethylene | 50 | 1.515338 | 30 | 1.510799 | 20 | 1.559923 | 10 | 1.618609 | 3 | 1.899324 | 0.5 | 1.60033 |
| 1,2,4-Trichlorobenzene | 50 | 0.6741326 | 30 | 0.6220869 | 20 | 0.5808433 | 10 | 0.5110483 | 3 | 0.3788416 | 0.5 | 0.2096018 |
| 1,2,4-Trimethylbenzene | 50 | 1.236641 | 30 | 1.254616 | 20 | 1.269096 | 10 | 1.319755 | 3 | 1.494853 | 0.5 | 1.421568 |
| 1,2-Dibromoethane | 50 | 0.4952987 | 30 | 0.497466 | 20 | 0.5125308 | 10 | 0.5279229 | 3 | 0.6240122 | 0.5 | 0.4963437 |
| 1,2-Dichlorobenzene | 50 | 0.7807969 | 30 | 0.7842246 | 20 | 0.7811683 | 10 | 0.8013964 | 3 | 0.886021 | 0.5 | 0.7791098 |
| 1,2-Dichloroethane | 50 | 1.472852 | 30 | 1.461965 | 20 | 1.510074 | 10 | 1.571183 | 3 | 1.870714 | 0.5 | 1.606795 |
| 1,2-Dichloropropane | 50 | 0.2406582 | 30 | 0.2419405 | 20 | 0.2484046 | 10 | 0.2586456 | 3 | 0.3068804 | 0.5 | 0.25533 |
| 1,2-Dichlorotetrafluoroethane | 50 | 2.430066 | 30 | 2.417112 | 20 | 2.503451 | 10 | 2.533954 | 3 | 2.879617 | 0.5 | 2.571998 |
| 1,3,5-Trimethylbenzene | 50 | 1.042181 | 30 | 1.04331 | 20 | 1.044135 | 10 | 1.191951 | 3 | 1.32858 | 0.5 | 1.216837 |
| 1,3-Butadiene | 50 | 0.529137 | 30 | 0.5264425 | 20 | 0.5460431 | 10 | 0.5517226 | 3 | 0.6445247 | 0.5 | 0.5742489 |
| 1,3-Dichlorobenzene | 50 | 0.8193318 | 30 | 0.8195923 | 20 | 0.8162876 | 10 | 0.8412337 | 3 | 0.9147794 | 0.5 | 0.7773592 |
| 1,3-Dichloropropane | 50 | 0.4463119 | 30 | 0.4453306 | 20 | 0.4593416 | 10 | 0.4769744 | 3 | 0.5594415 | 0.5 | 0.4676074 |
| 1,4-Dichlorobenzene | 50 | 0.8191321 | 30 | 0.8158773 | 20 | 0.8137969 | 10 | 0.8310733 | 3 | 0.8843848 | 0.5 | 0.740609 |
| 1,4-Dioxane | 50 | 0.1663113 | 30 | 0.1669162 | 20 | 0.1716396 | 10 | 0.1794077 | 3 | 0.2116021 | 0.5 | 0.2529028 |
| 2-Butanone | 50 | 1.510953 | 30 | 1.499307 | 20 | 1.534217 | 10 | 1.554827 | 3 | 1.784202 | 0.5 | 1.653068 |
| 2-Hexanone | 50 | 0.3892909 | 30 | 0.3858952 | 20 | 0.3914279 | 10 | 0.3957947 | 3 | 0.4356438 | 0.5 | 0.3680168 |
| 3-Chloropropene | 50 | 0.9332497 | 30 | 0.9225795 | 20 | 0.9534917 | 10 | 0.9827301 | 3 | 1.134481 | 0.5 | 0.9700514 |
| 4-Methyl-2-pentanone | 50 | 0.4124722 | 30 | 0.4093261 | 20 | 0.4175944 | 10 | 0.4221668 | 3 | 0.4947263 | 0.5 | 0.492995 |
| Acetone | 50 | 1.171301 | 30 | 1.164144 | 20 | 1.194752 | 10 | 1.221181 | 3 | 1.386964 | 0.5 | 4.449306 |
| Acrolein | 50 | 0.2751311 | 30 | 0.2696397 | 20 | 0.280614 | 10 | 0.2847119 | 3 | 0.3290155 | 0.5 | 0.2914644 |
| Acrylonitrile | 50 | 0.5998574 | 30 | 0.5966934 | 20 | 0.6093806 | 10 | 0.6278191 | 3 | 0.7423262 | 0.5 | 0.6415329 |
| Benzene | 50 | 3.335312 | 30 | 3.329007 | 20 | 3.443946 | 10 | 3.595121 | 3 | 4.343608 | 0.5 | 3.700083 |
| Benzyl chloride | 50 | 0.2110004 | 30 | 0.2079856 | 20 | 0.2025733 | 10 | 0.2006647 | 3 | 0.1977593 | 0.5 | 0.1188346 |
| Bromodichloromethane | 50 | 0.5670133 | 30 | 0.5691165 | 20 | 0.5851275 | 10 | 0.6050796 | 3 | 0.7108315 | 0.5 | 0.5641998 |
| Bromoform | 50 | 0.6791351 | 30 | 0.6857123 | 20 | 0.6914707 | 10 | 0.7232024 | 3 | 0.8268155 | 0.5 | 0.6275206 |

FORM VI

INITIAL CALIBRATION DATA

EPA TO-15

Laboratory: York Analytical Laboratories, Inc. - StratfordSDG: 20H0017Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YG00006Instrument: 5975CCalibration Date: 07/07/20 18:12

| Compound | Level 01 | | Level 02 | | Level 03 | | Level 04 | | Level 05 | | Level 06 | |
|--------------------------------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| Bromomethane | 50 | 0.9421696 | 30 | 0.9374589 | 20 | 0.961303 | 10 | 0.9766033 | 3 | 1.113665 | 0.5 | 1.009739 |
| Carbon disulfide | 50 | 2.769778 | 30 | 2.762278 | 20 | 2.856999 | 10 | 2.933426 | 3 | 3.434743 | 0.5 | 2.84724 |
| Carbon tetrachloride | 50 | 2.790028 | 30 | 2.784512 | 20 | 2.888584 | 10 | 2.98247 | 3 | 3.55798 | 0.5 | 2.887587 |
| Chlorobenzene | 50 | 0.7732371 | 30 | 0.7884924 | 20 | 0.8053137 | 10 | 0.8498696 | 3 | 0.9837685 | 0.5 | 0.8706818 |
| Chloroethane | 50 | 0.4650242 | 30 | 0.4607666 | 20 | 0.4771005 | 10 | 0.4853194 | 3 | 0.5495173 | 0.5 | 0.4933764 |
| Chloroform | 50 | 2.361867 | 30 | 2.355067 | 20 | 2.449932 | 10 | 2.531828 | 3 | 3.017948 | 0.5 | 2.486516 |
| Chloromethane | 50 | 0.5557616 | 30 | 0.5577326 | 20 | 0.5716246 | 10 | 0.5859443 | 3 | 0.6836927 | 0.5 | 0.7344974 |
| cis-1,2-Dichloroethylene | 50 | 1.351063 | 30 | 1.338536 | 20 | 1.389244 | 10 | 1.430402 | 3 | 1.699638 | 0.5 | 1.412067 |
| cis-1,3-Dichloropropylene | 50 | 0.457924 | 30 | 0.4581837 | 20 | 0.469965 | 10 | 0.4809509 | 3 | 0.5629795 | 0.5 | 0.4377289 |
| Cyclohexane | 50 | 1.356494 | 30 | 1.364906 | 20 | 1.418725 | 10 | 1.479746 | 3 | 1.771128 | 0.5 | 1.45828 |
| Dibromochloromethane | 50 | 0.6230053 | 30 | 0.6209389 | 20 | 0.6363368 | 10 | 0.6538475 | 3 | 0.7546436 | 0.5 | 0.5814883 |
| Dichlorodifluoromethane | 50 | 2.254814 | 30 | 2.402822 | 20 | 2.479072 | 10 | 2.637754 | 3 | 2.984151 | 0.5 | 2.456645 |
| Ethanol | 50 | 0.2510344 | 30 | 0.2492424 | 20 | 0.2555207 | 10 | 0.2679497 | 3 | 0.3109831 | 0.5 | 1.231346 |
| Ethyl acetate | 50 | 1.670447 | 30 | 1.646954 | 20 | 1.687181 | 10 | 1.719882 | 3 | 1.978908 | 0.5 | 1.632356 |
| Ethyl Benzene | 50 | 1.227427 | 30 | 1.273646 | 20 | 1.310957 | 10 | 1.385086 | 3 | 1.613945 | 0.5 | 1.465434 |
| Hexachlorobutadiene | 50 | 0.6546028 | 30 | 0.6633086 | 20 | 0.6750168 | 10 | 0.7033085 | 3 | 0.811371 | 0.5 | 0.7533861 |
| Isopropanol | 50 | 1.332178 | 30 | 1.318993 | 20 | 1.362162 | 10 | 1.401524 | 3 | 1.693545 | 0.5 | 4.940947 |
| Isopropylbenzene | 50 | 1.524866 | 30 | 1.550144 | 20 | 1.576984 | 10 | 1.65038 | 3 | 1.867188 | 0.5 | 1.656328 |
| Methyl Methacrylate | 50 | 0.2454148 | 30 | 0.2432515 | 20 | 0.2492317 | 10 | 0.2535235 | 3 | 0.2827617 | 0.5 | 0.2353807 |
| Methyl tert-butyl ether (MTBE) | 50 | 2.838307 | 30 | 2.84123 | 20 | 2.940289 | 10 | 3.036436 | 3 | 3.521152 | 0.5 | 3.025507 |
| Methylene chloride | 50 | 0.8840187 | 30 | 0.8822197 | 20 | 0.9092157 | 10 | 0.9356584 | 3 | 1.095488 | 0.5 | 1.167115 |
| Naphthalene | 50 | 1.396517 | 30 | 1.314499 | 20 | 1.229809 | 10 | 1.114901 | 3 | 0.8762422 | 0.5 | 1.100515 |
| n-Butylbenzene | 50 | 2.236237 | 30 | 2.299287 | 20 | 2.313363 | 10 | 2.361067 | 3 | 2.480136 | 0.5 | 1.801466 |
| n-Heptane | 50 | 1.413516 | 30 | 1.412063 | 20 | 1.466311 | 10 | 1.510924 | 3 | 1.774984 | 0.5 | 1.572794 |
| n-Hexane | 50 | 1.323331 | 30 | 1.33682 | 20 | 1.39155 | 10 | 1.446078 | 3 | 1.723145 | 0.5 | 1.432659 |
| n-Propylbenzene | 50 | 1.676272 | 30 | 1.713615 | 20 | 1.73729 | 10 | 1.813027 | 3 | 2.025828 | 0.5 | 1.814765 |
| o-Xylene | 50 | 1.036905 | 30 | 1.052358 | 20 | 1.074319 | 10 | 1.122768 | 3 | 1.272843 | 0.5 | 1.210526 |
| p- & m- Xylenes | 100 | 0.9802804 | 60 | 1.011487 | 40 | 1.037906 | 20 | 1.10091 | 6 | 1.262949 | 1 | 1.171171 |
| p-Ethyltoluene | 50 | 1.502661 | 30 | 1.55837 | 20 | 1.591674 | 10 | 1.681853 | 3 | 1.742718 | 0.5 | 1.518208 |
| p-Isopropyltoluene | 50 | 1.465324 | 30 | 1.491278 | 20 | 1.507742 | 10 | 1.557098 | 3 | 1.783846 | 0.5 | 1.523785 |
| Propylene | 50 | 0.3276534 | 30 | 0.3200154 | 20 | 0.2729133 | 10 | 0.3168536 | 3 | 0.3545432 | 0.5 | 0.3629387 |

FORM VI

INITIAL CALIBRATION DATA

EPA TO-15

Laboratory: York Analytical Laboratories, Inc. - StratfordSDG: 20H0017Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YG00006Instrument: 5975CCalibration Date: 07/07/20 18:12

| Compound | Level 01 | | Level 02 | | Level 03 | | Level 04 | | Level 05 | | Level 06 | |
|--------------------------------------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| sec-Butylbenzene | 50 | 1.730821 | 30 | 1.771843 | 20 | 1.796869 | 10 | 1.877809 | 3 | 2.144857 | 0.5 | 2.003641 |
| Styrene | 50 | 0.7864807 | 30 | 0.7958905 | 20 | 0.8025695 | 10 | 0.8324377 | 3 | 0.9121496 | 0.5 | 0.8295408 |
| tert-Butylbenzene | 50 | 1.210534 | 30 | 1.225204 | 20 | 1.23881 | 10 | 1.27891 | 3 | 1.463673 | 0.5 | 1.4008 |
| Tetrachloroethylene | 50 | 0.5004867 | 30 | 0.5002847 | 20 | 0.5145126 | 10 | 0.5328764 | 3 | 0.6240301 | 0.5 | 0.5714423 |
| Tetrahydrofuran | 50 | 0.8153016 | 30 | 0.8070331 | 20 | 0.8266924 | 10 | 0.8364987 | 3 | 0.9645189 | 0.5 | 0.8942671 |
| Toluene | 50 | 0.9416507 | 30 | 0.9469824 | 20 | 0.9720226 | 10 | 1.000377 | 3 | 1.163276 | 0.5 | 1.076391 |
| trans-1,2-Dichloroethylene | 50 | 1.361186 | 30 | 1.362835 | 20 | 1.414816 | 10 | 1.465575 | 3 | 1.741363 | 0.5 | 1.411468 |
| trans-1,3-Dichloropropylene | 50 | 0.4299573 | 30 | 0.4286999 | 20 | 0.4371284 | 10 | 0.4476828 | 3 | 0.5147422 | 0.5 | 0.3901596 |
| Trichloroethylene | 50 | 0.3434816 | 30 | 0.3424144 | 20 | 0.3517482 | 10 | 0.3649586 | 3 | 0.4305336 | 0.5 | 0.3554917 |
| Trichlorofluoromethane (Freon 11) | 50 | 2.872026 | 30 | 2.875758 | 20 | 2.978536 | 10 | 3.073591 | 3 | 3.590498 | 0.5 | 3.022873 |
| Vinyl acetate | 50 | 2.082996 | 30 | 2.036602 | 20 | 2.091108 | 10 | 2.14031 | 3 | 2.465286 | 0.5 | 2.036898 |
| Vinyl bromide | 50 | 1.020309 | 30 | 1.014272 | 20 | 1.047382 | 10 | 1.068416 | 3 | 1.216374 | 0.5 | 1.070678 |
| Vinyl Chloride | 50 | 0.7704521 | 30 | 0.7650928 | 20 | 0.7883654 | 10 | 0.7954742 | 3 | 0.9098342 | 0.5 | 0.8552376 |

FORM VI

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc. - Stratford

SDG: 20H0017

Client: Advanced Cleanup Technologies, Inc.

Project: 9628-BXNY

Calibration: YG00006

Instrument: 5975C

Calibration Date: 07/07/20 18:12

| Compound | Level 07 | | Level 08 | | Level 09 | | Level 10 | | Level 11 | | Level 12 | |
|---|----------|--------------|----------|--------------|----------|-----------|----------|-----------|----------|----|----------|----|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| 1,1,1,2-Tetrachloroethane | 0.2 | 0.4942459 | 0.1 | 0.504017 | 0.05 | 0.5242484 | 0.025 | 0.5320262 | | | | |
| 1,1,1-Trichloroethane | 0.2 | 2.64995 | 0.1 | 2.665339 | 0.05 | 2.812333 | 0.025 | 3.283566 | | | | |
| 1,1,2,2-Tetrachloroethane | 0.2 | 0.7023384 | 0.1 | 0.7261838 | 0.05 | 0.7372113 | 0.025 | 0.9156903 | | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 0.2 | 2.456361 | 0.1 | 2.488018 | 0.05 | 2.55796 | 0.025 | | | | | |
| 1,1,2-Trichloroethane | 0.2 | 0.3172114 | 0.1 | 0.3160495 | 0.05 | 0.379661 | 0.025 | 0.3287731 | | | | |
| 1,1-Dichloroethane | 0.2 | 1.963523 | 0.1 | 1.867834 | 0.05 | 2.102136 | 0.025 | 2.429426 | | | | |
| 1,1-Dichloroethylene | 0.2 | 1.590236 | 0.1 | 1.716718 | 0.05 | 1.538759 | 0.025 | 1.851592 | | | | |
| 1,2,4-Trichlorobenzene | 0.2 | 0.1531293 | 0.1 | | 0.05 | | 0.025 | | | | | |
| 1,2,4-Trimethylbenzene | 0.2 | 1.541574 | 0.1 | 1.799585 | 0.05 | 2.283958 | 0.025 | 2.866651 | | | | |
| 1,2-Dibromoethane | 0.2 | 0.4982678 | 0.1 | 0.5169821 | 0.05 | 0.5228567 | 0.025 | 0.6025802 | | | | |
| 1,2-Dichlorobenzene | 0.2 | 0.8081099 | 0.1 | 0.8750806 | 0.05 | 0.9406324 | 0.025 | 0.6907735 | | | | |
| 1,2-Dichloroethane | 0.2 | 1.658657 | 0.1 | 1.60986 | 0.05 | 1.827846 | 0.025 | 1.965095 | | | | |
| 1,2-Dichloropropane | 0.2 | 0.2487471 | 0.1 | 0.2845965 | 0.05 | 0.1964749 | 0.025 | 0.32954 | | | | |
| 1,2-Dichlorotetrafluoroethane | 0.2 | 2.615813 | 0.1 | 2.64059 | 0.05 | 2.796399 | 0.025 | 2.811209 | | | | |
| 1,3,5-Trimethylbenzene | 0.2 | 1.273272 | 0.1 | 1.530143 | 0.05 | 1.558087 | 0.025 | 1.894583 | | | | |
| 1,3-Butadiene | 0.2 | 0.6072186 | 0.1 | 0.655416 | 0.05 | 0.7062131 | 0.025 | | | | | |
| 1,3-Dichlorobenzene | 0.2 | 0.773463 | 0.1 | 0.8075641 | 0.05 | 0.8723183 | 0.025 | 0.6783111 | | | | |
| 1,3-Dichloropropane | 0.2 | 0.4726098 | 0.1 | 0.4648346 | 0.05 | 0.5023647 | 0.025 | 0.5067094 | | | | |
| 1,4-Dichlorobenzene | 0.2 | 0.7352247 | 0.1 | 0.7918287 | 0.05 | 0.8680314 | 0.025 | | | | | |
| 1,4-Dioxane | 0.2 | 0.3019326 | 0.1 | 0.338073 | 0.05 | 0.4695774 | 0.025 | 0.3499925 | | | | |
| 2-Butanone | 0.2 | 1.802445 | 0.1 | 1.961589 | 0.05 | 2.333178 | 0.025 | 2.521146 | | | | |
| 2-Hexanone | 0.2 | 0.3987625 | 0.1 | 0.4307235 | 0.05 | 0.4466016 | 0.025 | 0.4775646 | | | | |
| 3-Chloropropene | 0.2 | 0.9803574 | 0.1 | 1.078481 | 0.05 | 0.6703618 | 0.025 | 1.059363 | | | | |
| 4-Methyl-2-pentanone | 0.2 | 0.5447165 | 0.1 | 0.6286817 | 0.05 | 0.7137428 | 0.025 | 0.6601027 | | | | |
| Acetone | 0.2 | 6.520248 | 0.1 | 10.42493 | 0.05 | 17.79532 | 0.025 | 23.31859 | | | | |
| Acrolein | 0.2 | 0.2818121 | 0.1 | 0.2524414 | 0.05 | | 0.025 | | | | | |
| Acrylonitrile | 0.2 | 0.6274642 | 0.1 | 0.7086995 | 0.05 | 0.6521516 | 0.025 | | | | | |
| Benzene | 0.2 | 4.052804 | 0.1 | 4.377981 | 0.05 | 4.099564 | 0.025 | 3.637833 | | | | |
| Benzyl chloride | 0.2 | 9.637039E-02 | 0.1 | 9.087695E-02 | 0.05 | | 0.025 | | | | | |
| Bromodichloromethane | 0.2 | 0.5589315 | 0.1 | 0.5766607 | 0.05 | 0.582594 | 0.025 | 0.6634262 | | | | |
| Bromoform | 0.2 | 0.6296293 | 0.1 | 0.5973895 | 0.05 | 0.6214646 | 0.025 | 0.6771242 | | | | |

FORM VI

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc. - StratfordSDG: 20H0017Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YG00006Instrument: 5975CCalibration Date: 07/07/20 18:12

| Compound | Level 07 | | Level 08 | | Level 09 | | Level 10 | | Level 11 | | Level 12 | |
|--------------------------------|----------|-----------|----------|---------------------|----------|----------------------|----------|----------------------|----------|----|----------|----|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| Bromomethane | 0.2 | 1.004445 | 0.1 | 1.069164 | 0.05 | 1.047085 | 0.025 | | | | | |
| Carbon disulfide | 0.2 | 2.882257 | 0.1 | 2.873523 | 0.05 | 3.145237 | 0.025 | 3.171209 | | | | |
| Carbon tetrachloride | 0.2 | 2.838367 | 0.1 | 2.906134 | 0.05 | 2.961998 | 0.025 | 3.360381 | | | | |
| Chlorobenzene | 0.2 | 0.8926144 | 0.1 | 0.9544506 | 0.05 | 1.070623 | 0.025 | 1.186303 | | | | |
| Chloroethane | 0.2 | 0.4227921 | 0.1 | 0.1688766 | 0.05 | 0.3118491 | 0.025 | | | | | |
| Chloroform | 0.2 | 2.519462 | 0.1 | 2.596915 | 0.05 | 2.616574 | 0.025 | 2.812356 | | | | |
| Chloromethane | 0.2 | 0.6985453 | 0.1 | 0.4577139 | 0.05 | 0.6567042 | 0.025 | | | | | |
| cis-1,2-Dichloroethylene | 0.2 | 1.396943 | 0.1 | 1.380421 | 0.05 | 1.428929 | 0.025 | 1.498471 | | | | |
| cis-1,3-Dichloropropylene | 0.2 | 0.4585559 | 0.1 | 0.4618601 | 0.05 | 0.4848533 | 0.025 | | | | | |
| Cyclohexane | 0.2 | 1.462556 | 0.1 | 1.473303 | 0.05 | 1.645744 | 0.025 | | | | | |
| Dibromochloromethane | 0.2 | 0.5654749 | 0.1 | 0.5799515 | 0.05 | 0.614636 | 0.025 | 0.7112338 | | | | |
| Dichlorodifluoromethane | 0.2 | 2.746301 | 0.1 | 3.063073 | 0.05 | 3.414976 | 0.025 | 3.877451 | | | | |
| Ethanol | 0.2 | 1.632057 | 0.1 | 2.264402 | 0.05 | 3.260758 | 0.025 | 4.677705 | | | | |
| Ethyl acetate | 0.2 | 1.747472 | 0.1 | 1.892583 | 0.05 | 1.888736 | 0.025 | | | | | |
| Ethyl Benzene | 0.2 | 1.633473 | 0.1 | 1.914793 | 0.05 | 2.435107 | 0.025 | 3.33043 | | | | |
| Hexachlorobutadiene | 0.2 | 0.8569113 | 0.1 | 0.9357345 | 0.05 | 0.982257 | 0.025 | | | | | |
| Isopropanol | 0.2 | 6.122283 | 0.1 | 8.010574 | 0.05 | 11.0752 | 0.025 | 13.73273 | | | | |
| Isopropylbenzene | 0.2 | 1.702203 | 0.1 | 1.774145 | 0.05 | 1.991342 | 0.025 | 2.104367 | | | | |
| Methyl Methacrylate | 0.2 | 0.238948 | 0.1 | 0.2373852 | 0.05 | 0.2389492 | 0.025 | 0.2518207 | | | | |
| Methyl tert-butyl ether (MTBE) | 0.2 | 3.182541 | 0.1 | 3.358315 | 0.05 | 3.785439 | 0.025 | | | | | |
| Methylene chloride | 0.2 | | 0.1 | | 0.05 | | 0.025 | | | | | |
| Naphthalene | 0.2 | 0.8913117 | 0.1 | 0.7299963 | 0.05 | 0.4892616 | 0.025 | | | | | |
| n-Butylbenzene | 0.2 | 1.657169 | 0.1 | 1.608474 | 0.05 | 1.531121 | 0.025 | | | | | |
| n-Heptane | 0.2 | 1.608265 | 0.1 | 1.64218 | 0.05 | 1.745331 | 0.025 | 1.895159 | | | | |
| n-Hexane | 0.2 | 1.479994 | 0.1 | 1.444769 | 0.05 | 1.769232 | 0.025 | 1.650955 | | | | |
| n-Propylbenzene | 0.2 | 1.835896 | 0.1 | 2.063967 | 0.05 | 2.30802 | 0.025 | 2.237299 | | | | |
| o-Xylene | 0.2 | 1.36841 | 0.1 | 1.693597 | 0.05 | 2.209698 | 0.025 | 3.062192 | | | | |
| p- & m- Xylenes | 0.4 | 1.267674 | 0.2 | 1.492884 | 0.1 | 1.873313 | 0.05 | 2.569778 | | | | |
| p-Ethyltoluene | 0.2 | 1.655409 | 0.1 | 1.436216 | 0.05 | 2.021489 | 0.025 | 1.601123 | | | | |
| p-Isopropyltoluene | 0.2 | 1.602101 | 0.1 | 1.605909 | 0.05 | 1.616306 | 0.025 | 1.730494 | | | | |
| Propylene | 0.2 | 0.3701832 | 0.1 | 0.2268188 | 0.05 | | 0.025 | | | | | |

FORM VI

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc. - Stratford

SDG: 20H0017

Client: Advanced Cleanup Technologies, Inc.

Project: 9628-BXNY

Calibration: YG00006

Instrument: 5975C

Calibration Date: 07/07/20 18:12

| Compound | Level 07 | | Level 08 | | Level 09 | | Level 10 | | Level 11 | | Level 12 | |
|-----------------------------------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|----|----------|----|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| sec-Butylbenzene | 0.2 | 2.033215 | 0.1 | 2.43517 | 0.05 | 2.73284 | 0.025 | 2.715025 | | | | |
| Styrene | 0.2 | 0.8542001 | 0.1 | 0.9454392 | 0.05 | 1.155117 | 0.025 | 1.195501 | | | | |
| tert-Butylbenzene | 0.2 | 1.546398 | 0.1 | 1.76756 | 0.05 | 2.192442 | 0.025 | 2.013866 | | | | |
| Tetrachloroethylene | 0.2 | 0.6844493 | 0.1 | | 0.05 | | 0.025 | | | | | |
| Tetrahydrofuran | 0.2 | 0.9806529 | 0.1 | 1.054023 | 0.05 | 1.230894 | 0.025 | 1.031847 | | | | |
| Toluene | 0.2 | 1.122182 | 0.1 | 1.24382 | 0.05 | 1.545842 | 0.025 | 1.71008 | | | | |
| trans-1,2-Dichloroethylene | 0.2 | 1.408322 | 0.1 | 1.412158 | 0.05 | 1.431774 | 0.025 | 1.568407 | | | | |
| trans-1,3-Dichloropropylene | 0.2 | 0.3966995 | 0.1 | 0.3988274 | 0.05 | 0.421887 | 0.025 | | | | | |
| Trichloroethylene | 0.2 | 0.3649816 | 0.1 | 0.3809808 | 0.05 | 0.4071079 | 0.025 | 0.4220874 | | | | |
| Trichlorofluoromethane (Freon 11) | 0.2 | 3.093431 | 0.1 | 3.180121 | 0.05 | 3.14353 | 0.025 | | | | | |
| Vinyl acetate | 0.2 | 1.99426 | 0.1 | 2.088829 | 0.05 | 2.302448 | 0.025 | 2.282674 | | | | |
| Vinyl bromide | 0.2 | 1.081585 | 0.1 | 1.034515 | 0.05 | 1.024322 | 0.025 | | | | | |
| Vinyl Chloride | 0.2 | 0.9249407 | 0.1 | 1.056644 | 0.05 | 0.7500313 | 0.025 | 1.304713 | | | | |

FORM VI

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc. - StratfordSDG: 20H0017Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YG00006Instrument: 5975CCalibration Date: 07/07/20 18:12

| Compound | Mean RF | RF RSD | Mean RT | RT RSD | Linear r | Quad COD | LIMIT | Q |
|--|-----------|----------|----------|--------------|----------|-----------|-------|---|
| 1,1,1,2-Tetrachloroethane | 0.4845596 | 8.439021 | 21.46412 | 2.058193E-02 | | | 30 | |
| 1,1,1-Trichloroethane | 2.691275 | 8.231552 | 14.797 | 0.037001 | | | 30 | |
| 1,1,2,2-Tetrachloroethane | 0.7280765 | 9.172541 | 23.54813 | 1.433489E-02 | | | 30 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 2.421925 | 8.475933 | 10.08775 | 7.314307E-02 | | | 30 | |
| 1,1,2-Trichloroethane | 0.3240523 | 8.267195 | 19.2945 | 1.513009E-02 | | | 30 | |
| 1,1-Dichloroethane | 1.916883 | 8.002475 | 12.78125 | 4.068081E-02 | | | 30 | |
| 1,1-Dichloroethylene | 1.640163 | 8.411662 | 10.5504 | 6.741569E-02 | | | 30 | |
| 1,2,4-Trichlorobenzene | 0.4470977 | 45.80345 | 29.44514 | 1.774311E-02 | | 0.9995111 | 0.99 | |
| 1,2,4-Trimethylbenzene | 1.417211 | 13.55176 | 24.96 | 2.283221E-02 | | | 30 | |
| 1,2-Dibromoethane | 0.521103 | 8.297284 | 20.726 | 1.512118E-02 | | | 30 | |
| 1,2-Dichlorobenzene | 0.8119884 | 5.380939 | 26.58737 | 1.349903E-02 | | | 30 | |
| 1,2-Dichloroethane | 1.595262 | 8.235736 | 15.504 | 2.058922E-02 | | | 30 | |
| 1,2-Dichloropropane | 0.2606504 | 8.920483 | 16.847 | 2.381619E-02 | | | 30 | |
| 1,2-Dichlorotetrafluoroethane | 2.574075 | 5.715264 | 6.625 | 6.899894E-02 | | | 30 | |
| 1,3,5-Trimethylbenzene | 1.208801 | 14.13553 | 24.2055 | 1.851181E-02 | | | 30 | |
| 1,3-Butadiene | 0.5793442 | 8.760071 | 7.411125 | 0.1004589 | | | 30 | |
| 1,3-Dichlorobenzene | 0.8212014 | 5.362632 | 25.7715 | 1.919791E-02 | | | 30 | |
| 1,3-Dichloropropane | 0.4740565 | 7.663741 | 19.72825 | 3.026514E-02 | | | 30 | |
| 1,4-Dichlorobenzene | 0.8039908 | 6.049164 | 25.94175 | 1.146918E-02 | | | 30 | |
| 1,4-Dioxane | 0.2235982 | 29.94899 | 17.28363 | 9.965244E-02 | | | 30 | |
| 2-Butanone | 1.662576 | 10.20708 | 13.4755 | 6.629676E-02 | | | 30 | |
| 2-Hexanone | 0.3994444 | 5.709738 | 19.137 | 4.127823E-02 | | | 30 | |
| 3-Chloropropene | 0.9944277 | 7.427633 | 11.2475 | 4.823764E-02 | | | 30 | |
| 4-Methyl-2-pentanone | 0.4778349 | 16.48546 | 17.64325 | 0.0489066 | | | 30 | |
| Acetone | 3.441603 | 100.5348 | 10.32237 | 0.143961 | | 0.9998204 | 0.99 | |
| Acrolein | 0.2831038 | 7.75408 | 10.20112 | 0.1321792 | | | 30 | |
| Acrylonitrile | 0.6442217 | 8.25142 | 11.7275 | 0.1037619 | | | 30 | |
| Benzene | 3.772233 | 11.44092 | 15.57025 | 2.077412E-02 | | | 30 | |
| Benzyl chloride | 0.1657582 | 32.28868 | 26.124 | 5.672156E-03 | | 0.9999681 | 0.99 | |
| Bromodichloromethane | 0.5921201 | 8.46562 | 17.29138 | 1.369796E-02 | | | 30 | |
| Bromoform | 0.6826094 | 10.45426 | 23.46525 | 1.197565E-02 | | | 30 | |
| Bromomethane | 1.001818 | 6.209899 | 8.52475 | 4.318016E-02 | | | 30 | |

FORM VII

CONTINUING CALIBRATION CHECK

EPA TO-15

Laboratory: York Analytical Laboratories, Inc. - StratfordSDG: 20H0017Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYInstrument ID: 5975CCalibration: YG00006Lab File ID: TO285076.DCalibration Date: 07/07/20 18:12Sequence: Y0H1018Injection Date: 08/07/20Lab Sample ID: Y0H1018-CCV1Injection Time: 19:52

| COMPOUND | TYPE | CONC. (ppbv) | | RESPONSE FACTOR | | | % DIFF / DRIFT | |
|---|------|--------------|------|-----------------|-----------|---------|----------------|-----------|
| | | STD | CCV | ICAL | CCV | MIN (#) | CCV | LIMIT (#) |
| 1,1,1,2-Tetrachloroethane | A | 10.0 | 9.52 | 0.4845596 | 0.4615311 | | -4.8 | 30 |
| 1,1,1-Trichloroethane | A | 10.0 | 9.97 | 2.691275 | 2.683159 | | -0.3 | 30 |
| 1,1,2,2-Tetrachloroethane | A | 10.0 | 9.06 | 0.7280765 | 0.6599699 | | -9.4 | 30 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | A | 10.0 | 10.4 | 2.421925 | 2.528656 | | 4.4 | 30 |
| 1,1,2-Trichloroethane | A | 10.0 | 8.63 | 0.3240523 | 0.279614 | | -13.7 | 30 |
| 1,1-Dichloroethane | A | 10.0 | 9.99 | 1.916883 | 1.914678 | | -0.1 | 30 |
| 1,1-Dichloroethylene | A | 10.0 | 9.59 | 1.640163 | 1.572985 | | -4.1 | 30 |
| 1,2,4-Trichlorobenzene | A | 10.0 | 9.93 | 0.4470977 | 0.5425508 | | -0.7 | 30 |
| 1,2,4-Trimethylbenzene | A | 10.0 | 8.48 | 1.417211 | 1.20148 | | -15.2 | 30 |
| 1,2-Dibromoethane | A | 10.0 | 8.77 | 0.521103 | 0.4570431 | | -12.3 | 30 |
| 1,2-Dichlorobenzene | A | 10.0 | 10.0 | 0.8119884 | 0.8115886 | | -0.05 | 30 |
| 1,2-Dichloroethane | A | 10.0 | 9.18 | 1.595262 | 1.465028 | | -8.2 | 30 |
| 1,2-Dichloropropane | A | 10.0 | 8.08 | 0.2606504 | 0.2106568 | | -19.2 | 30 |
| 1,2-Dichlorotetrafluoroethane | A | 10.0 | 10.4 | 2.574075 | 2.675763 | | 4.0 | 30 |
| 1,3,5-Trimethylbenzene | A | 10.0 | 8.84 | 1.208801 | 1.068667 | | -11.6 | 30 |
| 1,3-Butadiene | A | 10.0 | 10.8 | 0.5793442 | 0.6270798 | | 8.2 | 30 |
| 1,3-Dichlorobenzene | A | 10.0 | 10.4 | 0.8212014 | 0.8532355 | | 3.9 | 30 |
| 1,3-Dichloropropane | A | 10.0 | 8.03 | 0.4740565 | 0.3806521 | | -19.7 | 30 |
| 1,4-Dichlorobenzene | A | 10.0 | 10.4 | 0.8039908 | 0.835535 | | 3.9 | 30 |
| 1,4-Dioxane | A | 10.0 | 6.71 | 0.2235982 | 0.1499355 | | -32.9 | 30 * |
| 2-Butanone | A | 10.0 | 9.40 | 1.662576 | 1.562618 | | -6.0 | 30 |
| 2-Hexanone | A | 10.0 | 7.51 | 0.3994444 | 0.2998667 | | -24.9 | 30 |
| 3-Chloropropene | A | 10.0 | 10.2 | 0.9944277 | 1.013555 | | 1.9 | 30 |
| 4-Methyl-2-pentanone | A | 10.0 | 6.89 | 0.4778349 | 0.3292578 | | -31.1 | 30 * |
| Acetone | Q | 10.0 | 9.46 | 3.441603 | 1.168801 | | -5.4 | 30 |
| Acrylonitrile | A | 10.0 | 10.0 | 0.6442217 | 0.6460466 | | 0.3 | 30 |
| Benzene | A | 10.0 | 10.0 | 3.772233 | 3.779549 | | 0.2 | 30 |
| Benzyl chloride | Q | 10.0 | 9.95 | 0.1657582 | 0.2000806 | | -0.5 | 30 |
| Bromodichloromethane | A | 10.0 | 8.02 | 0.5921201 | 0.4751251 | | -19.8 | 30 |

ATTACHMENT D

ANNOTATED SUMMARY FORMS

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 20H0017-01 File ID: TO285082.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 00:43
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|----|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 45.4 | 31 | UJ |
| 71-55-6 | 1,1,1-Trichloroethane | 45.4 | 25 | UJ |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 45.4 | 31 | UJ |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 45.4 | 35 | UJ |
| 79-00-5 | 1,1,2-Trichloroethane | 45.4 | 25 | UJ |
| 75-34-3 | 1,1-Dichloroethane | 45.4 | 18 | UJ |
| 75-35-4 | 1,1-Dichloroethylene | 45.4 | 4.5 | UJ |
| 120-82-1 | 1,2,4-Trichlorobenzene | 45.4 | 34 | UJ |
| 95-63-6 | 1,2,4-Trimethylbenzene | 45.4 | 22 | UJ |
| 106-93-4 | 1,2-Dibromoethane | 45.4 | 35 | UJ |
| 95-50-1 | 1,2-Dichlorobenzene | 45.4 | 27 | UJ |
| 107-06-2 | 1,2-Dichloroethane | 45.4 | 18 | UJ |
| 78-87-5 | 1,2-Dichloropropane | 45.4 | 21 | UJ |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 45.4 | 32 | UJ |
| 108-67-8 | 1,3,5-Trimethylbenzene | 45.4 | 22 | UJ |
| 106-99-0 | 1,3-Butadiene | 45.4 | 30 | UJ |
| 541-73-1 | 1,3-Dichlorobenzene | 45.4 | 27 | UJ |
| 142-28-9 | 1,3-Dichloropropane | 45.4 | 21 | UJ |
| 106-46-7 | 1,4-Dichlorobenzene | 45.4 | 27 | UJ |
| 123-91-1 | 1,4-Dioxane | 45.4 | 33 | UJ |
| 78-93-3 | 2-Butanone | 45.4 | 13 | UJ |
| 591-78-6 | 2-Hexanone | 45.4 | 37 | UJ |
| 107-05-1 | 3-Chloropropene | 45.4 | 71 | UJ |
| 108-10-1 | 4-Methyl-2-pentanone | 45.4 | 19 | UJ |
| 67-64-1 | Acetone | 45.4 | 620 | J |
| 107-13-1 | Acrylonitrile | 45.4 | 9.8 | UJ |
| 71-43-2 | Benzene | 45.4 | 35 | J |
| 100-44-7 | Benzyl chloride | 45.4 | 23 | UJ |
| 75-27-4 | Bromodichloromethane | 45.4 | 30 | UJ |
| 75-25-2 | Bromoform | 45.4 | 47 | UJ |
| 74-83-9 | Bromomethane | 45.4 | 18 | UJ |
| 75-15-0 | Carbon disulfide | 45.4 | 69 | J |
| 56-23-5 | Carbon tetrachloride | 45.4 | 7.1 | UJ |
| 108-90-7 | Chlorobenzene | 45.4 | 21 | UJ |
| 75-00-3 | Chloroethane | 45.4 | 12 | UJ |
| 67-66-3 | Chloroform | 45.4 | 22 | UJ |
| 74-87-3 | Chloromethane | 45.4 | 9.4 | UJ |
| 156-59-2 | cis-1,2-Dichloroethylene | 45.4 | 4.5 | UJ |
| 10061-01-5 | cis-1,3-Dichloropropylene | 45.4 | 21 | UJ |
| 110-82-7 | Cyclohexane | 45.4 | 7400 | J |

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 20H0017-01 File ID: TO285082.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 00:43
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 124-48-1 | Dibromochloromethane | 45.4 | 39 | UJ |
| 75-71-8 | Dichlorodifluoromethane | 45.4 | 22 | UJ |
| 141-78-6 | Ethyl acetate | 45.4 | 33 | UJ |
| 100-41-4 | Ethyl Benzene | 45.4 | 61 | J |
| 87-68-3 | Hexachlorobutadiene | 45.4 | 48 | UJ |
| 67-63-0 | Isopropanol | 45.4 | 22 | UJ |
| 80-62-6 | Methyl Methacrylate | 45.4 | 19 | UJ |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 45.4 | 16 | UJ |
| 75-09-2 | Methylene chloride | 45.4 | 32 | UJ |
| 142-82-5 | n-Heptane | 45.4 | 19 | UJ |
| 110-54-3 | n-Hexane | 45.4 | 1300 | J |
| 95-47-6 | o-Xylene | 45.4 | 20 | UJ |
| 179601-23-1 | p- & m- Xylenes | 45.4 | 39 | UJ |
| 622-96-8 | p-Ethyltoluene | 45.4 | 27 | J |
| 115-07-1 | Propylene | 45.4 | 370 | J |
| 100-42-5 | Styrene | 45.4 | 19 | UJ |
| 127-18-4 | Tetrachloroethylene | 45.4 | 130 | J |
| 109-99-9 | Tetrahydrofuran | 45.4 | 27 | UJ |
| 108-88-3 | Toluene | 45.4 | 56 | J |
| 156-60-5 | trans-1,2-Dichloroethylene | 45.4 | 31 | J |
| 10061-02-6 | trans-1,3-Dichloropropylene | 45.4 | 21 | UJ |
| 79-01-6 | Trichloroethylene | 45.4 | 6.1 | UJ |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 45.4 | 25 | UJ |
| 108-05-4 | Vinyl acetate | 45.4 | 16 | UJ |
| 593-60-2 | Vinyl bromide | 45.4 | 20 | UJ |
| 75-01-4 | Vinyl Chloride | 45.4 | 59 | J |

| CAS NO. | TENTATIVELY IDENTIFIED COMPOUND | RT | EST. CONC. (ppbv) | Q |
|----------|---------------------------------|--------|-------------------|---|
| 460-00-4 | SURR: p-Bromofluorobenzene | 23.753 | 13 | D |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 194625 | 14.462 | 240947 | 14.462 | |
| ISTD: 1,4-Difluorobenzene | 1099983 | 15.882 | 1394980 | 15.882 | |
| ISTD: d5-Chlorobenzene | 796732 | 21.363 | 1235217 | 21.363 | |

* Values outside of QC limits

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 20H0017-02 File ID: TO285083.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 01:29
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|----|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 21.7 | 15 | UJ |
| 71-55-6 | 1,1,1-Trichloroethane | 21.7 | 12 | UJ |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 21.7 | 15 | UJ |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 21.7 | 17 | UJ |
| 79-00-5 | 1,1,2-Trichloroethane | 21.7 | 12 | UJ |
| 75-34-3 | 1,1-Dichloroethane | 21.7 | 8.8 | UJ |
| 75-35-4 | 1,1-Dichloroethylene | 21.7 | 2.2 | UJ |
| 120-82-1 | 1,2,4-Trichlorobenzene | 21.7 | 16 | UJ |
| 95-63-6 | 1,2,4-Trimethylbenzene | 21.7 | 11 | UJ |
| 106-93-4 | 1,2-Dibromoethane | 21.7 | 17 | UJ |
| 95-50-1 | 1,2-Dichlorobenzene | 21.7 | 13 | UJ |
| 107-06-2 | 1,2-Dichloroethane | 21.7 | 8.8 | UJ |
| 78-87-5 | 1,2-Dichloropropane | 21.7 | 10 | UJ |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 21.7 | 15 | UJ |
| 108-67-8 | 1,3,5-Trimethylbenzene | 21.7 | 11 | UJ |
| 106-99-0 | 1,3-Butadiene | 21.7 | 14 | UJ |
| 541-73-1 | 1,3-Dichlorobenzene | 21.7 | 13 | UJ |
| 142-28-9 | 1,3-Dichloropropane | 21.7 | 10 | UJ |
| 106-46-7 | 1,4-Dichlorobenzene | 21.7 | 13 | UJ |
| 123-91-1 | 1,4-Dioxane | 21.7 | 16 | UJ |
| 78-93-3 | 2-Butanone | 21.7 | 6.4 | UJ |
| 591-78-6 | 2-Hexanone | 21.7 | 18 | UJ |
| 107-05-1 | 3-Chloropropene | 21.7 | 34 | UJ |
| 108-10-1 | 4-Methyl-2-pentanone | 21.7 | 8.9 | UJ |
| 67-64-1 | Acetone | 21.7 | 200 | J |
| 107-13-1 | Acrylonitrile | 21.7 | 4.7 | UJ |
| 71-43-2 | Benzene | 21.7 | 6.9 | UJ |
| 100-44-7 | Benzyl chloride | 21.7 | 11 | UJ |
| 75-27-4 | Bromodichloromethane | 21.7 | 15 | UJ |
| 75-25-2 | Bromoform | 21.7 | 22 | UJ |
| 74-83-9 | Bromomethane | 21.7 | 8.4 | UJ |
| 75-15-0 | Carbon disulfide | 21.7 | 22 | J |
| 56-23-5 | Carbon tetrachloride | 21.7 | 3.4 | UJ |
| 108-90-7 | Chlorobenzene | 21.7 | 10 | UJ |
| 75-00-3 | Chloroethane | 21.7 | 5.7 | UJ |
| 67-66-3 | Chloroform | 21.7 | 23 | J |
| 74-87-3 | Chloromethane | 21.7 | 4.5 | UJ |
| 156-59-2 | cis-1,2-Dichloroethylene | 21.7 | 1600 | J |
| 10061-01-5 | cis-1,3-Dichloropropylene | 21.7 | 9.9 | UJ |
| 110-82-7 | Cyclohexane | 21.7 | 51 | J |

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 20H0017-02 File ID: TO285083.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 01:29
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 124-48-1 | Dibromochloromethane | 21.7 | 19 | UJ |
| 75-71-8 | Dichlorodifluoromethane | 21.7 | 11 | UJ |
| 141-78-6 | Ethyl acetate | 21.7 | 16 | UJ |
| 100-41-4 | Ethyl Benzene | 21.7 | 9.4 | UJ |
| 87-68-3 | Hexachlorobutadiene | 21.7 | 23 | UJ |
| 67-63-0 | Isopropanol | 21.7 | 11 | UJ |
| 80-62-6 | Methyl Methacrylate | 21.7 | 8.9 | UJ |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 21.7 | 7.8 | UJ |
| 75-09-2 | Methylene chloride | 21.7 | 15 | UJ |
| 142-82-5 | n-Heptane | 21.7 | 8.9 | UJ |
| 110-54-3 | n-Hexane | 21.7 | 42 | J |
| 95-47-6 | o-Xylene | 21.7 | 9.4 | UJ |
| 179601-23-1 | p- & m- Xylenes | 21.7 | 19 | UJ |
| 622-96-8 | p-Ethyltoluene | 21.7 | 11 | UJ |
| 115-07-1 | Propylene | 21.7 | 32 | J |
| 100-42-5 | Styrene | 21.7 | 9.3 | UJ |
| 109-99-9 | Tetrahydrofuran | 21.7 | 13 | UJ |
| 108-88-3 | Toluene | 21.7 | 16 | J |
| 156-60-5 | trans-1,2-Dichloroethylene | 21.7 | 25 | J |
| 10061-02-6 | trans-1,3-Dichloropropylene | 21.7 | 9.9 | UJ |
| 79-01-6 | Trichloroethylene | 21.7 | 1500 | J |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 21.7 | 12 | UJ |
| 108-05-4 | Vinyl acetate | 21.7 | 7.6 | UJ |
| 593-60-2 | Vinyl bromide | 21.7 | 9.5 | UJ |
| 75-01-4 | Vinyl Chloride | 21.7 | 2.8 | UJ |

| CAS NO. | TENTATIVELY IDENTIFIED COMPOUND | RT | EST. CONC. (ppbv) | Q |
|----------|---------------------------------|--------|-------------------|---|
| 460-00-4 | SURR: p-Bromofluorobenzene | 23.753 | 10 | D |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 208096 | 14.462 | 240947 | 14.462 | |
| ISTD: 1,4-Difluorobenzene | 1176793 | 15.888 | 1394980 | 15.882 | |
| ISTD: d5-Chlorobenzene | 926359 | 21.363 | 1235217 | 21.363 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 20H0017-02RE1 File ID: TO285086.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 03:46
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|----------|---------------------|----------|----------------------------|---|
| 127-18-4 | Tetrachloroethylene | 86.9 | 9700 | D |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 205694 | 14.462 | 240947 | 14.462 | |
| ISTD: 1,4-Difluorobenzene | 1190206 | 15.888 | 1394980 | 15.882 | |
| ISTD: d5-Chlorobenzene | 946519 | 21.357 | 1235217 | 21.363 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Indoor Ambient Air Laboratory ID: 20H0017-03 File ID: TO285089.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 06:16
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|----|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.975 | 0.67 | UJ |
| 71-55-6 | 1,1,1-Trichloroethane | 0.975 | 0.53 | UJ |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.975 | 0.67 | UJ |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 0.975 | 0.97 | J |
| 79-00-5 | 1,1,2-Trichloroethane | 0.975 | 0.53 | UJ |
| 75-34-3 | 1,1-Dichloroethane | 0.975 | 0.39 | UJ |
| 75-35-4 | 1,1-Dichloroethylene | 0.975 | 0.097 | UJ |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.975 | 0.72 | UJ |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.975 | 2.3 | J |
| 106-93-4 | 1,2-Dibromoethane | 0.975 | 0.75 | UJ |
| 95-50-1 | 1,2-Dichlorobenzene | 0.975 | 0.59 | UJ |
| 107-06-2 | 1,2-Dichloroethane | 0.975 | 0.51 | J |
| 78-87-5 | 1,2-Dichloropropane | 0.975 | 0.45 | UJ |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 0.975 | 0.68 | UJ |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.975 | 0.77 | J |
| 106-99-0 | 1,3-Butadiene | 0.975 | 0.65 | UJ |
| 541-73-1 | 1,3-Dichlorobenzene | 0.975 | 0.59 | UJ |
| 142-28-9 | 1,3-Dichloropropane | 0.975 | 0.45 | UJ |
| 106-46-7 | 1,4-Dichlorobenzene | 0.975 | 1.2 | J |
| 123-91-1 | 1,4-Dioxane | 0.975 | 0.70 | UJ |
| 78-93-3 | 2-Butanone | 0.975 | 4.7 | J |
| 591-78-6 | 2-Hexanone | 0.975 | 0.80 | UJ |
| 107-05-1 | 3-Chloropropene | 0.975 | 1.5 | UJ |
| 108-10-1 | 4-Methyl-2-pentanone | 0.975 | 1.6 | J |
| 67-64-1 | Acetone | 0.975 | 98 | J |
| 107-13-1 | Acrylonitrile | 0.975 | 0.21 | UJ |
| 71-43-2 | Benzene | 0.975 | 1.3 | J |
| 100-44-7 | Benzyl chloride | 0.975 | 0.50 | UJ |
| 75-27-4 | Bromodichloromethane | 0.975 | 0.65 | UJ |
| 75-25-2 | Bromoform | 0.975 | 1.0 | UJ |
| 74-83-9 | Bromomethane | 0.975 | 0.38 | UJ |
| 75-15-0 | Carbon disulfide | 0.975 | 1.4 | J |
| 56-23-5 | Carbon tetrachloride | 0.975 | 0.86 | J |
| 108-90-7 | Chlorobenzene | 0.975 | 0.45 | UJ |
| 75-00-3 | Chloroethane | 0.975 | 0.26 | UJ |
| 67-66-3 | Chloroform | 0.975 | 0.57 | J |
| 74-87-3 | Chloromethane | 0.975 | 2.9 | J |
| 156-59-2 | cis-1,2-Dichloroethylene | 0.975 | 0.097 | UJ |
| 10061-01-5 | cis-1,3-Dichloropropylene | 0.975 | 0.44 | UJ |
| 110-82-7 | Cyclohexane | 0.975 | 0.44 | J |

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Indoor Ambient Air Laboratory ID: 20H0017-03 File ID: TO285089.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 06:16
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 124-48-1 | Dibromochloromethane | 0.975 | 0.83 | UJ |
| 75-71-8 | Dichlorodifluoromethane | 0.975 | 2.4 | J |
| 141-78-6 | Ethyl acetate | 0.975 | 4.1 | J |
| 100-41-4 | Ethyl Benzene | 0.975 | 2.8 | J |
| 87-68-3 | Hexachlorobutadiene | 0.975 | 1.0 | UJ |
| 67-63-0 | Isopropanol | 0.975 | 510 | J |
| 80-62-6 | Methyl Methacrylate | 0.975 | 0.60 | J |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 0.975 | 0.35 | UJ |
| 75-09-2 | Methylene chloride | 0.975 | 1.7 | J |
| 142-82-5 | n-Heptane | 0.975 | 1.1 | J |
| 110-54-3 | n-Hexane | 0.975 | 5.8 | J |
| 95-47-6 | o-Xylene | 0.975 | 3.1 | J |
| 179601-23-1 | p- & m- Xylenes | 0.975 | 10 | J |
| 622-96-8 | p-Ethyltoluene | 0.975 | 2.1 | J |
| 115-07-1 | Propylene | 0.975 | 0.17 | UJ |
| 100-42-5 | Styrene | 0.975 | 3.4 | J |
| 127-18-4 | Tetrachloroethylene | 0.975 | 12 | J |
| 109-99-9 | Tetrahydrofuran | 0.975 | 0.58 | UJ |
| 108-88-3 | Toluene | 0.975 | 16 | J |
| 156-60-5 | trans-1,2-Dichloroethylene | 0.975 | 0.39 | UJ |
| 10061-02-6 | trans-1,3-Dichloropropylene | 0.975 | 0.44 | UJ |
| 79-01-6 | Trichloroethylene | 0.975 | 0.42 | J |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 0.975 | 2.3 | J |
| 108-05-4 | Vinyl acetate | 0.975 | 0.34 | UJ |
| 593-60-2 | Vinyl bromide | 0.975 | 0.43 | UJ |
| 75-01-4 | Vinyl Chloride | 0.975 | 0.12 | UJ |

| CAS NO. | TENTATIVELY IDENTIFIED COMPOUND | RT | EST. CONC. (ppbv) | Q |
|----------|---------------------------------|--------|-------------------|---|
| 460-00-4 | SURR: p-Bromofluorobenzene | 23.753 | 9.8 | D |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 203360 | 14.462 | 240947 | 14.462 | |
| ISTD: 1,4-Difluorobenzene | 1109293 | 15.888 | 1394980 | 15.882 | |
| ISTD: d5-Chlorobenzene | 919003 | 21.363 | 1235217 | 21.363 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 20H0017-04 File ID: TO285084.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 02:15
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 124-48-1 | Dibromochloromethane | 16.9 | 14 | UJ |
| 75-71-8 | Dichlorodifluoromethane | 16.9 | 8.4 | UJ |
| 141-78-6 | Ethyl acetate | 16.9 | 12 | UJ |
| 100-41-4 | Ethyl Benzene | 16.9 | 7.3 | UJ |
| 87-68-3 | Hexachlorobutadiene | 16.9 | 18 | UJ |
| 67-63-0 | Isopropanol | 16.9 | 8.3 | UJ |
| 80-62-6 | Methyl Methacrylate | 16.9 | 6.9 | UJ |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 16.9 | 6.1 | UJ |
| 75-09-2 | Methylene chloride | 16.9 | 12 | UJ |
| 142-82-5 | n-Heptane | 16.9 | 44 | J |
| 110-54-3 | n-Hexane | 16.9 | 30 | J |
| 95-47-6 | o-Xylene | 16.9 | 7.3 | UJ |
| 179601-23-1 | p- & m- Xylenes | 16.9 | 15 | UJ |
| 622-96-8 | p-Ethyltoluene | 16.9 | 8.3 | UJ |
| 115-07-1 | Propylene | 16.9 | 21 | J |
| 100-42-5 | Styrene | 16.9 | 7.2 | UJ |
| 109-99-9 | Tetrahydrofuran | 16.9 | 10 | UJ |
| 108-88-3 | Toluene | 16.9 | 11 | J |
| 156-60-5 | trans-1,2-Dichloroethylene | 16.9 | 6.7 | UJ |
| 10061-02-6 | trans-1,3-Dichloropropylene | 16.9 | 7.7 | UJ |
| 79-01-6 | Trichloroethylene | 16.9 | 42 | J |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 16.9 | 9.5 | UJ |
| 108-05-4 | Vinyl acetate | 16.9 | 6.0 | UJ |
| 593-60-2 | Vinyl bromide | 16.9 | 7.4 | UJ |
| 75-01-4 | Vinyl Chloride | 16.9 | 2.2 | UJ |

| CAS NO. | TENTATIVELY IDENTIFIED COMPOUND | RT | EST. CONC. (ppbv) | Q |
|----------|---------------------------------|--------|-------------------|---|
| 460-00-4 | SURR: p-Bromofluorobenzene | 23.759 | 9.9 | |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 200629 | 14.462 | 240947 | 14.462 | |
| ISTD: 1,4-Difluorobenzene | 1127288 | 15.882 | 1394980 | 15.882 | |
| ISTD: d5-Chlorobenzene | 857998 | 21.363 | 1235217 | 21.363 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 20H0017-04RE1 File ID: TO285087.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 04:32
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|----------|---------------------|----------|----------------------------|---|
| 127-18-4 | Tetrachloroethylene | 84.5 | 15000 | D |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 190408 | 14.462 | 240947 | 14.462 | |
| ISTD: 1,4-Difluorobenzene | 1075600 | 15.888 | 1394980 | 15.882 | |
| ISTD: d5-Chlorobenzene | 790684 | 21.363 | 1235217 | 21.363 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 20H0017-05 File ID: TO285085.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 03:00
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|----|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 18.1 | 12 | UJ |
| 71-55-6 | 1,1,1-Trichloroethane | 18.1 | 9.8 | UJ |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 18.1 | 12 | UJ |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 18.1 | 14 | UJ |
| 79-00-5 | 1,1,2-Trichloroethane | 18.1 | 9.8 | UJ |
| 75-34-3 | 1,1-Dichloroethane | 18.1 | 7.3 | UJ |
| 75-35-4 | 1,1-Dichloroethylene | 18.1 | 1.8 | UJ |
| 120-82-1 | 1,2,4-Trichlorobenzene | 18.1 | 13 | UJ |
| 95-63-6 | 1,2,4-Trimethylbenzene | 18.1 | 8.9 | UJ |
| 106-93-4 | 1,2-Dibromoethane | 18.1 | 14 | UJ |
| 95-50-1 | 1,2-Dichlorobenzene | 18.1 | 11 | UJ |
| 107-06-2 | 1,2-Dichloroethane | 18.1 | 7.3 | UJ |
| 78-87-5 | 1,2-Dichloropropane | 18.1 | 8.3 | UJ |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 18.1 | 13 | UJ |
| 108-67-8 | 1,3,5-Trimethylbenzene | 18.1 | 8.9 | UJ |
| 106-99-0 | 1,3-Butadiene | 18.1 | 12 | UJ |
| 541-73-1 | 1,3-Dichlorobenzene | 18.1 | 11 | UJ |
| 142-28-9 | 1,3-Dichloropropane | 18.1 | 8.3 | UJ |
| 106-46-7 | 1,4-Dichlorobenzene | 18.1 | 11 | UJ |
| 123-91-1 | 1,4-Dioxane | 18.1 | 13 | UJ |
| 78-93-3 | 2-Butanone | 18.1 | 11 | J |
| 591-78-6 | 2-Hexanone | 18.1 | 15 | UJ |
| 107-05-1 | 3-Chloropropene | 18.1 | 28 | UJ |
| 108-10-1 | 4-Methyl-2-pentanone | 18.1 | 7.4 | UJ |
| 67-64-1 | Acetone | 18.1 | 93 | J |
| 107-13-1 | Acrylonitrile | 18.1 | 3.9 | UJ |
| 71-43-2 | Benzene | 18.1 | 12 | J |
| 100-44-7 | Benzyl chloride | 18.1 | 9.3 | UJ |
| 75-27-4 | Bromodichloromethane | 18.1 | 12 | UJ |
| 75-25-2 | Bromoform | 18.1 | 19 | UJ |
| 74-83-9 | Bromomethane | 18.1 | 7.0 | UJ |
| 75-15-0 | Carbon disulfide | 18.1 | 13 | J |
| 56-23-5 | Carbon tetrachloride | 18.1 | 2.8 | UJ |
| 108-90-7 | Chlorobenzene | 18.1 | 8.3 | UJ |
| 75-00-3 | Chloroethane | 18.1 | 4.8 | UJ |
| 67-66-3 | Chloroform | 18.1 | 8.8 | UJ |
| 74-87-3 | Chloromethane | 18.1 | 3.7 | UJ |
| 156-59-2 | cis-1,2-Dichloroethylene | 18.1 | 5.7 | J |
| 10061-01-5 | cis-1,3-Dichloropropylene | 18.1 | 8.2 | UJ |
| 110-82-7 | Cyclohexane | 18.1 | 6.2 | UJ |

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 20H0017-05 File ID: TO285085.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 03:00
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 124-48-1 | Dibromochloromethane | 18.1 | 15 | UJ |
| 75-71-8 | Dichlorodifluoromethane | 18.1 | 8.9 | UJ |
| 141-78-6 | Ethyl acetate | 18.1 | 13 | UJ |
| 100-41-4 | Ethyl Benzene | 18.1 | 7.8 | UJ |
| 87-68-3 | Hexachlorobutadiene | 18.1 | 19 | UJ |
| 67-63-0 | Isopropanol | 18.1 | 8.9 | UJ |
| 80-62-6 | Methyl Methacrylate | 18.1 | 7.4 | UJ |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 18.1 | 6.5 | UJ |
| 75-09-2 | Methylene chloride | 18.1 | 13 | UJ |
| 142-82-5 | n-Heptane | 18.1 | 21 | J |
| 110-54-3 | n-Hexane | 18.1 | 15 | J |
| 95-47-6 | o-Xylene | 18.1 | 7.8 | UJ |
| 179601-23-1 | p- & m- Xylenes | 18.1 | 16 | UJ |
| 622-96-8 | p-Ethyltoluene | 18.1 | 8.9 | UJ |
| 115-07-1 | Propylene | 18.1 | 13 | J |
| 100-42-5 | Styrene | 18.1 | 7.7 | UJ |
| 109-99-9 | Tetrahydrofuran | 18.1 | 11 | UJ |
| 108-88-3 | Toluene | 18.1 | 10 | J |
| 156-60-5 | trans-1,2-Dichloroethylene | 18.1 | 7.2 | UJ |
| 10061-02-6 | trans-1,3-Dichloropropylene | 18.1 | 8.2 | UJ |
| 79-01-6 | Trichloroethylene | 18.1 | 77 | J |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 18.1 | 10 | UJ |
| 108-05-4 | Vinyl acetate | 18.1 | 6.4 | UJ |
| 593-60-2 | Vinyl bromide | 18.1 | 7.9 | UJ |
| 75-01-4 | Vinyl Chloride | 18.1 | 2.3 | UJ |

| CAS NO. | TENTATIVELY IDENTIFIED COMPOUND | RT | EST. CONC. (ppbv) | Q |
|----------|---------------------------------|--------|-------------------|---|
| 460-00-4 | SURR: p-Bromofluorobenzene | 23.753 | 9.1 | |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 211971 | 14.462 | 240947 | 14.462 | |
| ISTD: 1,4-Difluorobenzene | 1231577 | 15.888 | 1394980 | 15.882 | |
| ISTD: d5-Chlorobenzene | 1001384 | 21.363 | 1235217 | 21.363 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 20H0017-05RE1 File ID: TO285088.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 05:17
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|----------|---------------------|----------|----------------------------|---|
| 127-18-4 | Tetrachloroethylene | 72.2 | 6400 | J |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 215848 | 14.462 | 240947 | 14.462 | |
| ISTD: 1,4-Difluorobenzene | 1237454 | 15.888 | 1394980 | 15.882 | |
| ISTD: d5-Chlorobenzene | 1033777 | 21.357 | 1235217 | 21.363 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Indoor Ambient Air Laboratory ID: 20H0017-06 File ID: TO285091.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 08:00
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|----|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.947 | 0.65 | UJ |
| 71-55-6 | 1,1,1-Trichloroethane | 0.947 | 0.52 | UJ |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.947 | 0.65 | UJ |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 0.947 | 0.73 | UJ |
| 79-00-5 | 1,1,2-Trichloroethane | 0.947 | 0.52 | UJ |
| 75-34-3 | 1,1-Dichloroethane | 0.947 | 0.38 | UJ |
| 75-35-4 | 1,1-Dichloroethylene | 0.947 | 0.094 | UJ |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.947 | 0.70 | UJ |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.947 | 1.2 | J |
| 106-93-4 | 1,2-Dibromoethane | 0.947 | 0.73 | UJ |
| 95-50-1 | 1,2-Dichlorobenzene | 0.947 | 0.57 | UJ |
| 107-06-2 | 1,2-Dichloroethane | 0.947 | 0.38 | UJ |
| 78-87-5 | 1,2-Dichloropropane | 0.947 | 0.44 | UJ |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 0.947 | 0.66 | UJ |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.947 | 0.47 | J |
| 106-99-0 | 1,3-Butadiene | 0.947 | 0.63 | UJ |
| 541-73-1 | 1,3-Dichlorobenzene | 0.947 | 0.57 | UJ |
| 142-28-9 | 1,3-Dichloropropane | 0.947 | 0.44 | UJ |
| 106-46-7 | 1,4-Dichlorobenzene | 0.947 | 0.57 | UJ |
| 123-91-1 | 1,4-Dioxane | 0.947 | 0.68 | UJ |
| 78-93-3 | 2-Butanone | 0.947 | 1.3 | J |
| 591-78-6 | 2-Hexanone | 0.947 | 0.78 | UJ |
| 107-05-1 | 3-Chloropropene | 0.947 | 1.5 | UJ |
| 108-10-1 | 4-Methyl-2-pentanone | 0.947 | 0.39 | UJ |
| 67-64-1 | Acetone | 0.947 | 10 | J |
| 107-13-1 | Acrylonitrile | 0.947 | 0.21 | UJ |
| 71-43-2 | Benzene | 0.947 | 0.48 | J |
| 100-44-7 | Benzyl chloride | 0.947 | 0.49 | UJ |
| 75-27-4 | Bromodichloromethane | 0.947 | 0.63 | UJ |
| 75-25-2 | Bromoform | 0.947 | 0.98 | UJ |
| 74-83-9 | Bromomethane | 0.947 | 0.37 | UJ |
| 75-15-0 | Carbon disulfide | 0.947 | 0.29 | UJ |
| 56-23-5 | Carbon tetrachloride | 0.947 | 0.48 | J |
| 108-90-7 | Chlorobenzene | 0.947 | 0.44 | UJ |
| 75-00-3 | Chloroethane | 0.947 | 0.25 | UJ |
| 67-66-3 | Chloroform | 0.947 | 0.46 | UJ |
| 74-87-3 | Chloromethane | 0.947 | 1.4 | J |
| 156-59-2 | cis-1,2-Dichloroethylene | 0.947 | 0.094 | UJ |
| 10061-01-5 | cis-1,3-Dichloropropylene | 0.947 | 0.43 | UJ |
| 110-82-7 | Cyclohexane | 0.947 | 0.33 | UJ |

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Indoor Ambient Air Laboratory ID: 20H0017-06 File ID: TO285091.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 08:00
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 124-48-1 | Dibromochloromethane | 0.947 | 0.81 | UJ |
| 75-71-8 | Dichlorodifluoromethane | 0.947 | 1.5 | J |
| 141-78-6 | Ethyl acetate | 0.947 | 0.68 | UJ |
| 100-41-4 | Ethyl Benzene | 0.947 | 0.41 | UJ |
| 87-68-3 | Hexachlorobutadiene | 0.947 | 1.0 | UJ |
| 67-63-0 | Isopropanol | 0.947 | 2.8 | J |
| 80-62-6 | Methyl Methacrylate | 0.947 | 0.58 | J |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 0.947 | 0.34 | UJ |
| 75-09-2 | Methylene chloride | 0.947 | 2.4 | J |
| 142-82-5 | n-Heptane | 0.947 | 0.47 | J |
| 110-54-3 | n-Hexane | 0.947 | 0.67 | J |
| 95-47-6 | o-Xylene | 0.947 | 0.45 | J |
| 179601-23-1 | p- & m- Xylenes | 0.947 | 1.1 | J |
| 622-96-8 | p-Ethyltoluene | 0.947 | 1.1 | J |
| 115-07-1 | Propylene | 0.947 | 0.16 | UJ |
| 100-42-5 | Styrene | 0.947 | 0.40 | UJ |
| 127-18-4 | Tetrachloroethylene | 0.947 | 16 | D |
| 109-99-9 | Tetrahydrofuran | 0.947 | 0.56 | UJ |
| 108-88-3 | Toluene | 0.947 | 1.4 | J |
| 156-60-5 | trans-1,2-Dichloroethylene | 0.947 | 0.38 | UJ |
| 10061-02-6 | trans-1,3-Dichloropropylene | 0.947 | 0.43 | UJ |
| 79-01-6 | Trichloroethylene | 0.947 | 26 | J |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 0.947 | 1.3 | J |
| 108-05-4 | Vinyl acetate | 0.947 | 0.33 | UJ |
| 593-60-2 | Vinyl bromide | 0.947 | 0.41 | UJ |
| 75-01-4 | Vinyl Chloride | 0.947 | 0.12 | UJ |

| CAS NO. | TENTATIVELY IDENTIFIED COMPOUND | RT | EST. CONC. (ppbv) | Q |
|----------|---------------------------------|--------|-------------------|---|
| 460-00-4 | SURR: p-Bromofluorobenzene | 23.753 | 8.8 | |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 224875 | 14.462 | 240947 | 14.462 | |
| ISTD: 1,4-Difluorobenzene | 1310612 | 15.882 | 1394980 | 15.882 | |
| ISTD: d5-Chlorobenzene | 1099973 | 21.363 | 1235217 | 21.363 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Outdoor Ambient Air Laboratory ID: 20H0017-07 File ID: TO285093.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 09:58
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|----|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.936 | 0.64 | UJ |
| 71-55-6 | 1,1,1-Trichloroethane | 0.936 | 0.51 | UJ |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 0.936 | 0.64 | UJ |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 0.936 | 0.72 | UJ |
| 79-00-5 | 1,1,2-Trichloroethane | 0.936 | 0.51 | UJ |
| 75-34-3 | 1,1-Dichloroethane | 0.936 | 0.38 | UJ |
| 75-35-4 | 1,1-Dichloroethylene | 0.936 | 0.093 | UJ |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.936 | 0.69 | UJ |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.936 | 0.60 | J |
| 106-93-4 | 1,2-Dibromoethane | 0.936 | 0.72 | UJ |
| 95-50-1 | 1,2-Dichlorobenzene | 0.936 | 0.56 | UJ |
| 107-06-2 | 1,2-Dichloroethane | 0.936 | 0.38 | UJ |
| 78-87-5 | 1,2-Dichloropropane | 0.936 | 0.43 | UJ |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 0.936 | 0.65 | UJ |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.936 | 0.46 | UJ |
| 106-99-0 | 1,3-Butadiene | 0.936 | 0.62 | UJ |
| 541-73-1 | 1,3-Dichlorobenzene | 0.936 | 0.56 | UJ |
| 142-28-9 | 1,3-Dichloropropane | 0.936 | 0.43 | UJ |
| 106-46-7 | 1,4-Dichlorobenzene | 0.936 | 0.56 | UJ |
| 123-91-1 | 1,4-Dioxane | 0.936 | 0.67 | UJ |
| 78-93-3 | 2-Butanone | 0.936 | 1.2 | J |
| 591-78-6 | 2-Hexanone | 0.936 | 0.77 | UJ |
| 107-05-1 | 3-Chloropropene | 0.936 | 1.5 | UJ |
| 108-10-1 | 4-Methyl-2-pentanone | 0.936 | 0.38 | UJ |
| 67-64-1 | Acetone | 0.936 | 8.8 | J |
| 107-13-1 | Acrylonitrile | 0.936 | 0.20 | UJ |
| 71-43-2 | Benzene | 0.936 | 0.69 | J |
| 100-44-7 | Benzyl chloride | 0.936 | 0.48 | UJ |
| 75-27-4 | Bromodichloromethane | 0.936 | 0.63 | UJ |
| 75-25-2 | Bromoform | 0.936 | 0.97 | UJ |
| 74-83-9 | Bromomethane | 0.936 | 0.36 | UJ |
| 75-15-0 | Carbon disulfide | 0.936 | 0.29 | UJ |
| 56-23-5 | Carbon tetrachloride | 0.936 | 0.53 | J |
| 108-90-7 | Chlorobenzene | 0.936 | 0.43 | UJ |
| 75-00-3 | Chloroethane | 0.936 | 0.25 | UJ |
| 67-66-3 | Chloroform | 0.936 | 0.73 | J |
| 74-87-3 | Chloromethane | 0.936 | 1.4 | J |
| 156-59-2 | cis-1,2-Dichloroethylene | 0.936 | 0.093 | UJ |
| 10061-01-5 | cis-1,3-Dichloropropylene | 0.936 | 0.42 | UJ |
| 110-82-7 | Cyclohexane | 0.936 | 0.35 | J |

Laboratory: York Analytical Laboratories, Inc. - Stratford SDG: 20H0017
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Outdoor Ambient Air Laboratory ID: 20H0017-07 File ID: TO285093.D
 Sampled: 07/31/20 14:00 Prepared: 08/06/20 09:00 Analyzed: 08/08/20 09:58
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BH00312 Sequence: Y0H1018 Calibration: YG00006 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 124-48-1 | Dibromochloromethane | 0.936 | 0.80 | UJ |
| 75-71-8 | Dichlorodifluoromethane | 0.936 | 1.7 | J |
| 141-78-6 | Ethyl acetate | 0.936 | 0.67 | UJ |
| 100-41-4 | Ethyl Benzene | 0.936 | 0.41 | J |
| 87-68-3 | Hexachlorobutadiene | 0.936 | 1.0 | UJ |
| 67-63-0 | Isopropanol | 0.936 | 6.4 | J |
| 80-62-6 | Methyl Methacrylate | 0.936 | 0.38 | UJ |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 0.936 | 0.34 | UJ |
| 75-09-2 | Methylene chloride | 0.936 | 1.3 | J |
| 142-82-5 | n-Heptane | 0.936 | 0.54 | J |
| 110-54-3 | n-Hexane | 0.936 | 0.99 | J |
| 95-47-6 | o-Xylene | 0.936 | 0.49 | J |
| 179601-23-1 | p- & m- Xylenes | 0.936 | 1.2 | J |
| 622-96-8 | p-Ethyltoluene | 0.936 | 0.55 | J |
| 115-07-1 | Propylene | 0.936 | 0.16 | UJ |
| 100-42-5 | Styrene | 0.936 | 0.40 | UJ |
| 127-18-4 | Tetrachloroethylene | 0.936 | 0.63 | J |
| 109-99-9 | Tetrahydrofuran | 0.936 | 0.55 | UJ |
| 108-88-3 | Toluene | 0.936 | 1.8 | J |
| 156-60-5 | trans-1,2-Dichloroethylene | 0.936 | 0.37 | UJ |
| 10061-02-6 | trans-1,3-Dichloropropylene | 0.936 | 0.42 | UJ |
| 79-01-6 | Trichloroethylene | 0.936 | 0.25 | J |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 0.936 | 1.3 | J |
| 108-05-4 | Vinyl acetate | 0.936 | 0.33 | UJ |
| 593-60-2 | Vinyl bromide | 0.936 | 0.41 | UJ |
| 75-01-4 | Vinyl Chloride | 0.936 | 0.12 | UJ |

| CAS NO. | TENTATIVELY IDENTIFIED COMPOUND | RT | EST. CONC. (ppbv) | Q |
|----------|---------------------------------|--------|-------------------|---|
| 460-00-4 | SURR: p-Bromofluorobenzene | 23.753 | 9.1 | |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 215317 | 14.462 | 240947 | 14.462 | |
| ISTD: 1,4-Difluorobenzene | 1232590 | 15.888 | 1394980 | 15.882 | |
| ISTD: d5-Chlorobenzene | 1005017 | 21.363 | 1235217 | 21.363 | |

* Values outside of QC limits

JR 12/8/2025

December 8, 2025

Mr. Jason Stewart
Advanced Cleanup Technologies
228 Park Ave S PMB 34864
New York, New York 10003

Re: Data Usability Summary Report – York Analytical Laboratories, Inc. - 21A0950

Dear Mr. Stewart:

The evaluation of analytical data by York Analytical Laboratories for project 9268-BXNY, which were reported in a single data package under Job No. 21A0950 has been completed. The following samples were reported.

SS-5 SS-6 IA-3

Analysis was performed in accordance with EPA Method TO-15 (volatile organics). The review was performed to the extent possible, in accordance with the analytical method, and “DER-10/ Technical Guidance for Site Investigation and Remediation”. Professional judgment is applied as necessary and appropriate. National Functional Guidelines for Organic Data Review was consulted as needed. Qualifiers consistent with those defined by EPA Region 2 are applied as necessary and appropriate.

Below is the Data Usability Summary Report (DUSR) associated with these samples.

| Data Usability Summary Report | |
|---|---|
| 1. Is the data package complete as defined under the requirements for the most current DEC ASP Category B or USEPA CLP data deliverables? | No -see Documentation section and section A.3 |
| 2. Have all holding times been met? | Yes |
| 3. Do all the QC data; blanks, instrument tunings, calibration standards, calibration verifications, surrogate recoveries, spike recoveries, replicate analyses, laboratory controls and sample data fall within the protocol required limits and specifications? | No -see following sections |
| 4. Have all of the data been generated using established and agreed upon analytical protocols? | Yes |
| 5. Does an evaluation of the raw data confirm the results provided in the data summary sheet and the quality control verification forms? | Yes |
| 6. Have the correct data qualifiers been used and are they consistent with the most current DEC ASP? | Yes |

| | |
|---|-----|
| 7. Have any quality control (QC) exceedances been specifically noted in the DUSR and have the corresponding QC summary sheet from the data package been attached to the DUSR? | Yes |
|---|-----|

Overall Evaluation

Based on the data review effort, results are usable, with the following qualifications. For samples that are qualified as estimated (J-, UJ), detected results may be biased low. False negatives may exist in non-detect results. Sample results that are qualified as estimated (J+) may be biased high. For samples that are qualified as estimated with any combination of (J), (J-) and/or (J+), the (J) qualifier takes precedence and is applied to the sample result. It is not possible to determine the direction of the bias and the overall effect on the result. Where a sample result is rejected (R), the 'R' qualifier takes precedence over any other qualification.

- The results for all target analytes in all samples are qualified as estimated (J, UJ) because there is no documentation provided to show samples were properly preserved.
- Based on professional judgment, the non-detect result for 1,2,4-trichlorobenzene in all samples is rejected (R) due to unacceptable RSD in the initial calibration (IC).
- Results for methylene chloride and p-ethyltoluene are qualified as estimated (J-), due to unacceptable calibration.
- Quantitation limits (QLs) for several analytes are adjusted based on unacceptable initial calibration. See Section A.1.
- The results for dichlorodifluoromethane and propylene in IA-3 are qualified as estimated (UJ) due to low response in the continuing calibration verification (CCV) standard.
- The results for dichlorodifluoromethane and isopropanol in IA- 3 are qualified as estimated (J-, UJ) due to low recovery in the laboratory control sample (LCS). The non-detect result for propylene in IA-3 is rejected (R) because the LCS recovery is less than 10%.
- Based on professional judgment, samples results less than ten times the quantitation limit in all samples are qualified as estimated (J) because proper canister certification cannot be confirmed.

Qualifier definitions are provided in Attachment A. A copy of the chain of custody record is provided in Attachment B. Pages from the data package illustrating the exceedances and issues described in this validation report are included in Attachment C. Annotated Summary

Forms are included in Attachment D detailing qualifications resulting from the data review effort.

The following components were reviewed, where applicable:

- Chain of Custody
- Receiving conditions
- Holding times
- Preservation
- Analyte lists
- Reporting limits
- Requested methods
- Units, and
- Sample related quality control data:
 - Method, instrument blanks
 - Clean canister certification
 - Field blanks
 - Surrogate recoveries
 - LCS recoveries
 - Internal standard area response
 - Duplicates
 - Analyte Identification
- Instrument related quality control data:
 - Instrument tunes
 - Calibration summaries

The following sections of the report detail only quality control exceedances that impacted results. Where a quality control item exceeded control limits but there is no impact to the samples results, these are not detailed in the report.

Documentation: A completeness review of the data package was performed, and the data package was determined to be a complete Category B data package, with following exception:

- The data package does not include the clean canister certification data. The laboratory was contacted and was requested to provide the missing documentation. The laboratory stated *'our standard archival period is three years and due to this data being older than that we won't be able to obtain the additional documents.'* See Section A.3.

The following documentation issues were observed during the review:

- The data package indicates that the laboratory did not hold certification for ethyl acetate, p-ethyltoluene, propylene, tetrahydrofuran, 1,1,1,2-tetrachloroethane, 1,3-dichloropropane, and 2-hexanone. It is noted that NY ELAP does not offer certification for ethyl acetate, p-ethyltoluene, propylene, tetrahydrofuran and 2-hexanone, but

does offer certification for 1,1,1,2-tetrachloroethane and 1,3-dichloropropane. The laboratory was contacted to request clarification. The laboratory stated *'Given the age of the project, it is difficult to ascertain the exact circumstances in which these compounds were requested. We don't typically report non-certified compounds, except at the request of the client, so it can be safely assumed that there was correspondence regarding the accreditation status of the compounds at the time the samples were taken. Just to note- while these compounds are indeed non-certified under our scope of accreditation between 2020-2022, we ran the compounds in the exact same way we ran our certified compounds – quantitated against a multiple calibration, where they were properly qualified as specified by our certifying body.'*

Holding Times, Preservation, Sample Integrity:

A copy of the applicable chain of custody (COC) record was included in the data package, documenting sample collection date of January 22, 2021. The samples were received at the laboratory on the same day on January 25, 2021. The samples were analyzed within method hold time. However, the data package does contain sample receipt information regarding condition of canisters upon receipt. The laboratory was contacted for clarification and stated *'Any anomalous receipt data would be denoted in the case narrative. If there was nothing listed there wasn't any anomalous activity, but we've opened a corrective action to improve this process.'* Since there is no evidence that the samples were properly preserved (pressure inside the canister maintained within +/- 5 psi from sampling to check in the laboratory or analysis), the results for all target analytes in all samples are qualified as estimated (J, UJ).

A. Volatile Organics

1. Calibration

Two initial calibrations (ICs) were performed in support of the sample analysis. For sample IA-3, analyzed on instrument Air2, all relative response factors (RRFs) and relative standard deviations (RSDs) or correlation coefficients are acceptable with the exceptions noted below.

| Analyte | RSD |
|----------------------------------|------|
| Instrument Air2 1/26/2021 | |
| 1,2,4-Trichlorobenzene | 76 |
| 1,3-Dichlorobenzene | 34.1 |
| 1,4-Dichlorobenzene | 44.9 |
| Benzyl chloride | 57.6 |

The laboratory used the average response factor for all analytes in the initial calibration. The method requires an RSD of $\leq 30\%$, with an allowance of up to two analytes to have an RSD $\leq 40\%$. A review of the individual calibration levels indicates a lack of instrument sensitivity in the lower calibration standards as presented below.

| Analyte | RRF 50 | RRF 30 | RRF 20 | RRF 10 | RRF 3 | RRF 0.5 | RRF 0.2 | RRF 0.1 |
|------------------------|----------|----------|----------|----------|-----------|-----------------|-----------------|-----------------|
| 1,2,4-Trichlorobenzene | 0.936863 | 0.728804 | 0.69116 | 0.514499 | 0.2059795 | 0.009348 | 0.045876 | - |
| 1,3-Dichlorobenzene | 1.246324 | 0.997939 | 1.078101 | 1.048843 | 0.9167642 | 0.613606 | 0.498674 | 0.467932 |
| 1,4-Dichlorobenzene | 1.250273 | 0.999835 | 1.067838 | 1.01241 | 0.7902146 | 0.467800 | 0.382028 | 0.321293 |
| Benzyl chloride | 1.681651 | 1.313488 | 1.339341 | 1.169678 | 0.7215757 | 0.310940 | 0.223463 | - |

The validator calculated the RSD with these standards excluded (**bolded above**), and the RSDs are:

- 1,2,4-Trichlorobenzene – 44.5 RSD
- 1,3-Dichlorobenzene -11.5 RSD
- 1,4-Dichlorobenzene -16.1 RSD
- Benzyl chloride – 27.9 RSD

Based on professional judgment, the non-detect result for 1,2,4-trichlorobenzene in IA-3 is rejected (R) due to unacceptable RSD (>40%RSD) in the IC after the elimination of the standards with excessively low response.

1,3-Dichlorobenzene, 1,4-dichlorobenzene, and benzyl chloride are not detected in IA-3. The quantitation limits (QLs) have been adjusted to reflect the lowest acceptable concentration standard (3ppbv) in the IC as presented below:

| Analyte | IA-3 adjusted QL (ppbv) |
|---------------------|-------------------------|
| 1,3-Dichlorobenzene | 17.3 |
| 1,4-Dichlorobenzene | 17.3 |
| Benzyl chloride | 15.5 |

For samples SS-5 and SS-6, analyzed on instrument 5975C, all relative response factors (RRFs) and relative standard deviations (RSDs) or correlation coefficients are acceptable with the exceptions noted below.

| Analyte | RSD |
|--|------|
| <i>Instrument 5975C 1/28/2021</i> | |
| 1,2,4-Trichlorobenzene | 41.1 |
| 1,2,4-Trimethylbenzene | 53.1 |
| 1,2-Dichlorobenzene | 44.6 |
| 1,3,5-Trimethylbenzene | 44.0 |

| Analyte | RSD |
|---------------------|------|
| 1,3-Dichlorobenzene | 49.8 |
| 1,4-Dichlorobenzene | 54.6 |
| 2-Butanone | 44.9 |
| Ethyl acetate | 44.8 |
| Hexachlorobutadiene | 47.8 |
| Isopropanol | 30.2 |
| Methyl Methacrylate | 54.3 |
| Methylene chloride | 71.3 |
| p-Ethyltoluene | 48.8 |
| Styrene | 35.8 |
| Tetrahydrofuran | 37.5 |
| Vinyl acetate | 57.8 |
| Vinyl Chloride | 32.5 |

It is noted that the internal standard area responses for 0.1 ppbv and 50 ppbv standards are biased high, outside the 60-140%R control limit as presented below:

| Standard | Bromochloromethane %R | 1,4-Difluorobenzene %R | Chlorobenzene-d5 %R |
|------------------------------|-----------------------|------------------------|---------------------|
| Cal Standard (Y1A2808-CALA) | 143 | 160 | 154 |
| Cal Standard (Y1A2808-CAL3) | 148 | 159 | A |

A=acceptable

The laboratory used the average response factor for all analytes in the initial calibration. A review of the individual calibration levels is presented below for the analytes with RSD >30.

| Analyte | RRF 50 | RRF 30 | RRF 20 | RRF 10 | RRF 3 | RRF 0.5 | RRF 0.2 | RRF 0.1 |
|------------------------|----------|----------|----------|----------|-----------|-----------------|-----------------|-----------------|
| 1,2,4-Trichlorobenzene | 0.700693 | 0.677484 | 0.590836 | 0.408506 | 0.1966694 | - | - | - |
| 1,2,4-Trimethylbenzene | 1.375326 | 1.458105 | 1.446346 | 1.435829 | 1.542911 | 0.273764 | 0.342346 | 0.539647 |
| 1,2-Dichlorobenzene | 0.900839 | 0.959441 | 0.966115 | 0.943664 | 0.911476 | 0.245998 | 0.270073 | 0.507492 |
| 1,3,5-Trimethylbenzene | 1.322004 | 1.366753 | 1.372565 | 1.360943 | 1.540981 | 0.392061 | 0.464303 | 0.678614 |
| 1,3-Dichlorobenzene | 0.909889 | 0.957549 | 0.939073 | 0.903935 | 0.8244748 | 0.223418 | 0.222262 | 0.376262 |
| 1,4-Dichlorobenzene | 0.902312 | 0.937458 | 0.909971 | 0.844290 | 0.7197025 | 0.159711 | 0.181004 | 0.334693 |
| 2-Butanone | 2.294415 | 2.268154 | 2.279497 | 2.213659 | 2.426894 | 0.80472 | 0.755105 | 0.829266 |
| Ethyl acetate | 2.556619 | 2.562783 | 2.534473 | 2.451693 | 2.710518 | 0.855485 | 0.825364 | 0.999244 |

| | | | | | | | | |
|---------------------|----------|----------|----------|----------|-----------|------------------|-----------------------|-----------------|
| Isopropanol | 2.070019 | 2.067289 | 2.03248 | 2.031593 | 2.140223 | 1.009817 | 1.037407 | 1.211657 |
| Methyl Methacrylate | 0.294606 | 0.332228 | 0.341239 | 0.330449 | 0.3445318 | 0.0804109 | 0.077656 | 0.088947 |
| Methylene chloride | 1.250825 | 1.259062 | 1.277804 | 1.296804 | 1.555752 | 2.868581 | 5.248707 | - |
| p-Ethyltoluene | 1.61259 | 1.700738 | 1.712024 | 1.702453 | 1.767332 | 0.434838 | 0.477236 | 0.651678 |
| Styrene | 0.838986 | 0.932732 | 0.977283 | 0.969475 | 0.9446457 | 0.369036 | 0.397238 | 0.52576 |
| Tetrahydrofuran | 1.300876 | 1.284318 | 1.301304 | 1.271495 | 1.408516 | 0.527207 | 0.533769 | 0.651464 |
| Vinyl acetate | 3.065108 | 2.969136 | 2.840886 | 2.621604 | 2.886531 | 0.359552 | 0.532220 | 1.034093 |
| Vinyl Chloride | 0.589450 | 0.620361 | 0.666787 | 0.686514 | 0.7183967 | 0.780540 | 0.809031 | 1.411269 |
| | | | | | | | ORRF 0.025 | RRF0.05 |
| | | | | | | | 1.229984 | 1.017383 |

The method requires a minimum of five (5) standards for the IC, therefore, eliminating the lowest standard is not an option that can be used to achieve $RSD \leq 30$. Therefore, based on professional judgment, the non-detect results for 1,2,4-trichlorobenzene in SS-5 and SS-6 are rejected (R) due to unacceptable RSD ($>40\%RSD$) in the IC.

The high RSDs for the remaining analytes are caused by low response factors in the three lowest concentration standards (**bolded above**). The validator calculated the RSD with these three standards excluded, and the RSDs are presented below.

| Analyte | RSD |
|------------------------|-----|
| 1,2,4-Trimethylbenzene | 4.1 |
| 1,2-Dichlorobenzene | 3.1 |
| 1,3,5-Trimethylbenzene | 6.1 |
| 1,3-Dichlorobenzene | 5.6 |
| 1,4-Dichlorobenzene | 10 |
| 2-Butanone | 3.4 |
| Ethyl acetate | 3.6 |
| Isopropanol | 2.1 |
| Methyl Methacrylate | 6.1 |
| Methylene chloride | 9.7 |
| p-Ethyltoluene | 3.3 |
| Styrene | 5.9 |

| Analyte | RSD |
|-----------------|-----|
| Tetrahydrofuran | 4.2 |
| Vinyl acetate | 5.8 |

The table below shows the reported QLs and the adjusted QLs based on elimination of the low standards in the IC for the non-detect analytes in SS-5 and SS-6. These adjusted QLs are reflected in the annotated summary forms in Appendix D.

| Sample | Analyte | Reported QL (µg/m3) | Adjust QL (µg/m3) |
|--------|------------------------|---------------------|-------------------|
| SS-5 | 1,2,4-Trimethylbenzene | 17 U | 341 U |
| | 1,2-Dichlorobenzene | 14 U | 417 U |
| | 1,3,5-Trimethylbenzene | 11 U | 341 U |
| | 1,3-Dichlorobenzene | 14 U | 417 U |
| | 1,4-Dichlorobenzene | 14 U | 417 U |
| | 2-Butanone | 6.8 U | 204 U |
| | Ethyl acetate | 17 U | 250 U |
| | Isopropanol | 11 U | 254 U |
| | Methyl Methacrylate | 9.5 U | 240 U |
| | Methylene chloride | 16 U | 241 U |
| | p-Ethyltoluene | 11 U | 341 U |
| | Styrene | 9.8 U | 295 U |
| | Tetrahydrofuran | 14 U | 204 U |
| | Vinyl acetate | 8.1 U | 244 U |
| SS-6 | 1,2-Dichlorobenzene | 1.0 U | 30.5 U |
| | 1,3,5-Trimethylbenzene | 0.83 U | 24.9 U |
| | 1,3-Dichlorobenzene | 1.0 U | 30.5 U |
| | 1,4-Dichlorobenzene | 1.0 U | 30.5 U |
| | Ethyl acetate | 1.2 U | 18.3 U |
| | Methyl Methacrylate | 0.69 U | 20.8 U |
| | Styrene | 072 U | 21.6 U |
| | Vinyl acetate | 0.59 U | 17.8 U |

For detected analytes, the recalculated sample concentrations using the updated average relative response factors are presented below:

| Sample | Analyte | Reported Conc (on-column ppbv) | Adjusted Conc (on-column ppbv) | %D | Adjust QL (µg/m3) |
|--------|------------------------|--------------------------------|--------------------------------|-----|-------------------|
| SS-6 | 1,2,4-Trimethylbenzene | 0.17 | 0.175 | 2.9 | - |
| | 2-Butanone | 3.21 | 3.14 | 2.2 | - |
| | Methylene chloride | 0.98 | 1.88 | 92 | 17.6 |

| | | | | | |
|--|-----------------|------|-------|-----|------|
| | p-Ethyltoluene | 0.17 | 0.18 | 5.9 | 13.1 |
| | Isopropanol | 0.40 | 0.395 | 1.2 | - |
| | Tetrahydrofuran | 0.56 | 0.55 | 1.8 | - |
| | Vinyl chloride | 0.44 | 0.43 | 2.3 | - |

Where the difference in concentration is less than 5%, no action is taken. Where the difference in concentration exceeds 5% sample results are qualified as estimated (J-), biased low due to unacceptable calibration, as follows:

- Methylene chloride and p-ethyltoluene in SS-6.

For vinyl chloride, the high RSD is caused by high responses in the three lowest calibration standards. The concentration of vinyl chloride in SS-5 is greater than ten times the 0.1 ppbv calibration standard and the high response has no impact on the sample result.

A second source ICV standard was analyzed after the IC, and all percent differences are acceptable ($\leq 30\%D$). CCVs were analyzed at the appropriate frequency and are acceptable ($\%D \leq 30$) with the exceptions noted below.

| Analyte | %D | Associated Sample | Qualifier Applied |
|---|-------|-------------------|-------------------|
| <i>CCV Instrument Air2 01/26/21 07:48</i> | | | |
| Dichlorodifluoromethane | -68.5 | IA-3 | UJ |
| Propylene | -96.3 | | |

The percent differences represent a decrease in instrument sensitivity. The results for dichlorodifluoromethane and propylene in IA-3 are qualified as estimated (UJ) due to low response in the CCV.

2. Laboratory Control Sample (LCS) /LCS Duplicate (LCSD)

Three LCSs were prepared and analyzed with the sample. All recoveries are acceptable (70-130%R), with the following exceptions:

| Analyte | LCS %R | Affected Sample | Qualifier Applied |
|-------------------------|--------|-----------------|-------------------|
| <i>BA11425-BS1</i> | | | |
| Dichlorodifluoromethane | 30.4 | IA-3 | J- |
| Isopropanol | 51.0 | | |
| Propylene | 5.2 | | R |

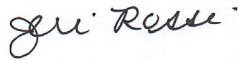
The results for dichlorodifluoromethane and isopropanol are qualified as estimated (J-, UJ) due to low recovery in the LCS. The non-detect result for propylene in IA-3 is rejected (R) because the LCS recovery is less than 10%.

3. Compound Quantitation

As noted in **Documentation** section, the laboratory did not provide clean canister certification data. Since clean canister certification data is not available, it cannot be confirmed whether the canisters used for sampling were properly cleaned and free of background contamination. This uncertainty introduces potential impacts to sample results, including the possibility that detected compounds may originate from residual contamination within the canisters rather than the sampled environment. As a result, accuracy cannot be fully assured, and reported concentrations may not reliably represent actual ambient conditions. Based on professional judgment, samples results less than ten times the quantitation limit in all samples are qualified as estimated (J) because proper canister certification cannot be confirmed.

No other sample results are qualified. Please feel free to contact me at (908) 370-3431 or richjერიrossi513@gmail.com if you have any questions regarding this data package review report or need further information.

Sincerely,

A handwritten signature in cursive script that reads "Jeri Rossi".

Jeri L Rossi, CEAC

Environmental Consulting Chemist

ATTACHMENT A

Qualifier Definitions

EPA Qualifier Definitions

- U The analyte was analyzed for but was not detected above the level of the reported sample quantitation limit.
- J The result is an estimated quantity. The associated numerical value is the approximate concentration of the analyte in the sample.
- J+ The result is an estimated quantity, but the result may be biased high.
- J- The result is an estimated quantity, but the result may be biased low.
- NJ The analyte has been “tentatively identified” or “presumptively” as present and the associated numerical value is the estimated concentration in the sample.
- UJ The analyte was analyzed for but was not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
- R The data are unusable. The sample results are rejected due to serious deficiencies in meeting QC criteria. The analyte may or may not be present in the sample.

ATTACHMENT B

CHAIN OF CUSTODY (COC)



York Analytical Laboratories, Inc.
 120 Research Drive Stratford, CT 06615
 132-02 89th Ave Queens, NY 11418
 clientservices@yorklab.com
 www.yorklab.com

Field Chain-of-Custody Record - AIR

YORK Project No.
 21A0950

NOTE: YORK's Standard Terms & Conditions are listed on the back side of this document. This document serves as your written authorization for YORK to proceed with the analyses requested below. Your signature binds you to YORK's Standard Terms & Conditions.

Your Page 1 of 1

| | | | | | | | | | |
|---|--|---------------------------------|--|---------------------------------|--|----------------------------|--|--|--|
| YOUR Information | | Report To: | | Invoice To: | | YOUR Project Number | | Turn-Around Time | |
| Company: <u>Advanced Clean Tech</u> | | Company: <u>ACT</u> | | Company: <u>ACT</u> | | <u>9628-BXNY</u> | | RUSH - Next Day | |
| Address: <u>200 Brainhollow Rd St. 207 Melville, NY</u> | | Address: <u>same</u> | | Address: <u>same</u> | | | | RUSH - Two Day | |
| Phone: <u>516-441-5600</u> | | Phone: _____ | | Phone: _____ | | YOUR Project Name | | RUSH - Three Day | |
| Contact: <u>Jason Stewart</u> | | Contact: <u>Paul Stewart</u> | | Contact: <u>Karen Freeman</u> | | | | RUSH - Four Day | |
| E-mail: <u>jason@act-earth</u> | | E-mail: <u>Paul's@act-earth</u> | | E-mail: <u>KarenF@act-earth</u> | | YOUR PO#: | | Standard (5-7 Day) <input checked="" type="checkbox"/> | |

Please print clearly and legibly. All information must be complete. Samples will not be logged in and the turn-around-time clock will not begin until any questions by YORK are resolved.

| | | | | | | |
|---|--|---|--|--|--|---|
| Samples Collected by: (print your name above and sign below) <u>Tom Yung</u> <u>[Signature]</u> | Air Matrix Codes | Samples From | Report / EDD Type (circle selections) | | | YORK Reg. Comp. |
| | AI - Indoor Ambient Air | New York | <input checked="" type="checkbox"/> Summary Report | CT RCP | <input checked="" type="checkbox"/> Standard Excel EDD | Compared to the following Regulation(s): (please fill in) |
| | AO - Outdoor Amb. Air | New Jersey | <input type="checkbox"/> QA Report | CT RCP DQA/DUE | <input type="checkbox"/> EQUIS (Standard) | |
| | AE - Vapor Extraction Well/ Process Gas/Effluent | Connecticut | <input type="checkbox"/> NY ASP A Package | NJDEP Reduced Deliv. | <input type="checkbox"/> NYSDEC EQUIS | |
| AS - Soil Vapor/Sub-Slab | Pennsylvania | <input type="checkbox"/> NY ASP B Package | NJDQOP | <input type="checkbox"/> NJDEP SRP HazSite | | |
| | | Other | <input type="checkbox"/> Other: | | | |

| Certified Canisters: Batch <input checked="" type="checkbox"/> Individual <input type="checkbox"/> | | Please enter the following REQUIRED Field Data | | | | | Reporting Units: ug/m ³ ___ ppbv ___ ppmv ___ | |
|--|-------------------|--|---|--|-------------|---------------|--|--|
| Sample Identification | Date/Time Sampled | Air Matrix | Canister Vacuum Before Sampling (in Hg) | Canister Vacuum After Sampling (in Hg) | Canister ID | Flow Cont. ID | Analysis Requested | |
| SS-5 | 11/22/21 8-1430 | AS | -30 | -12 | 36995 | 5705 | TO-15 | |
| SS-6 | " 10-1430 | AS | -29 | -8 | 36985 | 12011 | " | |
| IA-3 | " 8-1600 | AI | -29 | -10 | 36983 | 13573 | " | |
| | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |
| | | | | | | | | |

Comments: canister 23995 (stopped) may have taken in water - tanks filled w/ water (TV)

| | |
|--|--|
| Detection Limits Required | Sampling Media |
| ≤ 1 ug/m ³ ___ NYSDEC V1 Limits ___ Routine Survey ___ Other ___ | 6 Liter Canister <input checked="" type="checkbox"/> Tedlar Bag |

| | | | | | |
|---|-------------------------------------|--|-------------------------------------|---|-----------------------------------|
| Samples Relinquished by / Company <u>T-Y to ACT</u> | Date/Time <u>11/25/21 1610AM</u> | Samples Received by / Company <u>K. Smith with</u> | Date/Time <u>11/25/21 1010AM</u> | Samples Relinquished by / Company <u>K. Smith with</u> | Date/Time <u>11/25/21 1708</u> |
| Samples Received by / Company <u>TC Hahl / YORK</u> | Date/Time <u>11/25/21 1708</u> | Samples Relinquished by / Company <u>TC Hahl / YORK</u> | Date/Time <u>11/25/21 1830</u> | Samples Received by / Company <u>Ed [Signature]</u> | Date/Time <u>11/25/21 1830</u> |
| Samples Relinquished by / Company <u>[Signature]</u> | Date/Time <u>11/25/21 2020</u> | Samples Received by / Company | Date/Time | Samples Received in LAB by <u>J. [Signature] YORK</u> | Date/Time <u>11/26/21 9:50</u> |

Page 30 of 332

ATTACHMENT C

**SELECTED PAGES FROM DATA PACKAGE –
QC EXCEEDANCES AND VALIDATION ISSUES**

FORM III

LCS / LCS DUPLICATE RECOVERY

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYMatrix: AirBatch: BA11425Laboratory ID: BA11425-BS1Preparation: EPA TO15 PREPInitial/Final: 400 mL / 400 mL

| COMPOUND | SPIKE ADDED (ppbv) | LCS CONCENTRATION (ppbv) | LCS % REC. # | QC LIMITS REC. |
|--------------------------------|--------------------|--------------------------|--------------|----------------|
| Bromomethane | 10.0 | 9.22 | 92.2 | 70 - 130 |
| Carbon disulfide | 10.0 | 10.8 | 108 | 70 - 130 |
| Carbon tetrachloride | 10.0 | 11.9 | 119 | 70 - 130 |
| Chlorobenzene | 10.0 | 10.2 | 102 | 70 - 130 |
| Chloroethane | 10.0 | 9.03 | 90.3 | 70 - 130 |
| Chloroform | 10.0 | 10.9 | 109 | 70 - 130 |
| Chloromethane | 10.0 | 10.8 | 108 | 70 - 130 |
| cis-1,2-Dichloroethylene | 10.0 | 10.3 | 103 | 70 - 130 |
| cis-1,3-Dichloropropylene | 10.0 | 10.8 | 108 | 70 - 130 |
| Cyclohexane | 10.0 | 10.4 | 104 | 70 - 130 |
| Dibromochloromethane | 10.0 | 11.2 | 112 | 70 - 130 |
| Dichlorodifluoromethane | 10.0 | 3.04 | 30.4 * | 70 - 130 |
| Ethyl acetate | 10.0 | 10.5 | 105 | 70 - 130 |
| Ethyl Benzene | 10.0 | 10.7 | 107 | 70 - 130 |
| Hexachlorobutadiene | 10.0 | 11.9 | 119 | 70 - 130 |
| Isopropanol | 10.0 | 5.10 | 51.0 * | 70 - 130 |
| Methyl Methacrylate | 10.0 | 10.5 | 105 | 70 - 130 |
| Methyl tert-butyl ether (MTBE) | 10.0 | 10.9 | 109 | 70 - 130 |
| Methylene chloride | 10.0 | 9.58 | 95.8 | 70 - 130 |
| n-Heptane | 10.0 | 10.7 | 107 | 70 - 130 |
| n-Hexane | 10.0 | 9.16 | 91.6 | 70 - 130 |
| o-Xylene | 10.0 | 10.8 | 108 | 70 - 130 |
| p- & m- Xylenes | 20.0 | 21.8 | 109 | 70 - 130 |
| p-Ethyltoluene | 10.0 | 11.9 | 119 | 70 - 130 |
| Propylene | 10.0 | 0.520 | 5.20 * | 70 - 130 |
| Styrene | 10.0 | 11.5 | 115 | 70 - 130 |
| Tetrachloroethylene | 10.0 | 10.8 | 108 | 70 - 130 |
| Tetrahydrofuran | 10.0 | 10.3 | 103 | 70 - 130 |
| Toluene | 10.0 | 10.3 | 103 | 70 - 130 |
| trans-1,2-Dichloroethylene | 10.0 | 10.6 | 106 | 70 - 130 |

FORM VIII

INTERNAL STANDARD AREA AND RT SUMMARY
EPA TO-15

Laboratory: York Analytical Laboratories, Inc.

SDG: 21A0950

Client: Advanced Cleanup Technologies, Inc.

Project: 9628-BXNY

Sequence: Y1A2808

Instrument: 5975C

Calibration: YB10001

| Internal Standard | Response | RT | Reference Response | Reference RT | Area % | Area % Limits | RT Diff | RT Diff Limit | Q |
|------------------------------------|----------|--------|--------------------|--------------|--------------------------|---------------|---------|---------------|---|
| Cal Standard (Y1A2808-CAL1) | | | | | | | | | |
| Lab File ID: TO287788.D | | | | | Analyzed: 01/28/21 01:00 | | | | |
| Bromochloromethane | 325858 | 12.304 | 285825 | 12.304 | 114 | 60 - 140 | 0.0000 | +/-0.33 | |
| ISTD: 1,4-Difluorobenzene | 1076070 | 13.852 | 934692 | 13.852 | 115 | 60 - 140 | 0.0000 | +/-0.33 | |
| ISTD: d5-Chlorobenzene | 769231 | 19.095 | 818110 | 19.095 | 94 | 60 - 140 | 0.0000 | +/-0.33 | |
| Cal Standard (Y1A2808-CAL2) | | | | | | | | | |
| Lab File ID: TO287789.D | | | | | Analyzed: 01/28/21 01:47 | | | | |
| Bromochloromethane | 311387 | 12.304 | 285825 | 12.304 | 109 | 60 - 140 | 0.0000 | +/-0.33 | |
| ISTD: 1,4-Difluorobenzene | 1013162 | 13.852 | 934692 | 13.852 | 108 | 60 - 140 | 0.0000 | +/-0.33 | |
| ISTD: d5-Chlorobenzene | 743026 | 19.095 | 818110 | 19.095 | 91 | 60 - 140 | 0.0000 | +/-0.33 | |
| Cal Standard (Y1A2808-CAL6) | | | | | | | | | |
| Lab File ID: TO287793.D | | | | | Analyzed: 01/28/21 05:02 | | | | |
| Bromochloromethane | 273484 | 12.304 | 285825 | 12.304 | 96 | 60 - 140 | 0.0000 | +/-0.33 | |
| ISTD: 1,4-Difluorobenzene | 854667 | 13.852 | 934692 | 13.852 | 91 | 60 - 140 | 0.0000 | +/-0.33 | |
| ISTD: d5-Chlorobenzene | 708075 | 19.095 | 818110 | 19.095 | 87 | 60 - 140 | 0.0000 | +/-0.33 | |
| Cal Standard (Y1A2808-CAL7) | | | | | | | | | |
| Lab File ID: TO287794.D | | | | | Analyzed: 01/28/21 05:50 | | | | |
| Bromochloromethane | 285825 | 12.304 | 285825 | 12.304 | 100 | 60 - 140 | 0.0000 | +/-0.33 | |
| ISTD: 1,4-Difluorobenzene | 934692 | 13.852 | 934692 | 13.852 | 100 | 60 - 140 | 0.0000 | +/-0.33 | |
| ISTD: d5-Chlorobenzene | 818110 | 19.095 | 818110 | 19.095 | 100 | 60 - 140 | 0.0000 | +/-0.33 | |
| Cal Standard (Y1A2808-CAL8) | | | | | | | | | |
| Lab File ID: TO287795.D | | | | | Analyzed: 01/28/21 06:40 | | | | |
| Bromochloromethane | 319639 | 12.304 | 285825 | 12.304 | 112 | 60 - 140 | 0.0000 | +/-0.33 | |
| ISTD: 1,4-Difluorobenzene | 1098511 | 13.852 | 934692 | 13.852 | 118 | 60 - 140 | 0.0000 | +/-0.33 | |
| ISTD: d5-Chlorobenzene | 966685 | 19.095 | 818110 | 19.095 | 118 | 60 - 140 | 0.0000 | +/-0.33 | |
| Cal Standard (Y1A2808-CAL9) | | | | | | | | | |
| Lab File ID: TO287796.D | | | | | Analyzed: 01/28/21 07:33 | | | | |
| Bromochloromethane | 368307 | 12.304 | 285825 | 12.304 | 129 | 60 - 140 | 0.0000 | +/-0.33 | |
| ISTD: 1,4-Difluorobenzene | 1289667 | 13.852 | 934692 | 13.852 | 138 | 60 - 140 | 0.0000 | +/-0.33 | |
| ISTD: d5-Chlorobenzene | 1112418 | 19.095 | 818110 | 19.095 | 136 | 60 - 140 | 0.0000 | +/-0.33 | |
| Cal Standard (Y1A2808-CALA) | | | | | | | | | |
| Lab File ID: TO287797.D | | | | | Analyzed: 01/28/21 08:30 | | | | |
| Bromochloromethane | 409794 | 12.304 | 285825 | 12.304 | 143 | 60 - 140 | 0.0000 | +/-0.33 | * |
| ISTD: 1,4-Difluorobenzene | 1494630 | 13.852 | 934692 | 13.852 | 160 | 60 - 140 | 0.0000 | +/-0.33 | * |
| ISTD: d5-Chlorobenzene | 1258981 | 19.095 | 818110 | 19.095 | 154 | 60 - 140 | 0.0000 | +/-0.33 | * |
| Cal Standard (Y1A2808-CAL3) | | | | | | | | | |
| Lab File ID: TO287799.D | | | | | Analyzed: 01/28/21 10:10 | | | | |
| Bromochloromethane | 421819 | 12.304 | 285825 | 12.304 | 148 | 60 - 140 | 0.0000 | +/-0.33 | * |
| ISTD: 1,4-Difluorobenzene | 1490759 | 13.852 | 934692 | 13.852 | 159 | 60 - 140 | 0.0000 | +/-0.33 | * |
| ISTD: d5-Chlorobenzene | 1072923 | 19.095 | 818110 | 19.095 | 131 | 60 - 140 | 0.0000 | +/-0.33 | |
| Cal Standard (Y1A2808-CAL4) | | | | | | | | | |
| Lab File ID: TO287804.D | | | | | Analyzed: 01/28/21 14:21 | | | | |
| Bromochloromethane | 355116 | 12.31 | 285825 | 12.304 | 124 | 60 - 140 | 0.0060 | +/-0.33 | |
| ISTD: 1,4-Difluorobenzene | 1174406 | 13.858 | 934692 | 13.852 | 126 | 60 - 140 | 0.0060 | +/-0.33 | |
| ISTD: d5-Chlorobenzene | 854399 | 19.095 | 818110 | 19.095 | 104 | 60 - 140 | 0.0000 | +/-0.33 | |

FORM VI

INITIAL CALIBRATION DATA

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.

SDG: 21A0950

Client: Advanced Cleanup Technologies, Inc.

Project: 9628-BXNY

Calibration: YA10030

Instrument: TO15 AIR2

Calibration Date: 01/26/21 07:48

| Compound | Level 01 | | Level 02 | | Level 03 | | Level 04 | | Level 05 | | Level 06 | |
|---|----------|---------|----------|----------|----------|-----------|----------|--------------|----------|--------------|----------|-----------|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| 1,1,1,2-Tetrachloroethane | 0.025 | | 0.05 | | 0.1 | 0.5770163 | 0.2 | 0.5752757 | 0.5 | 0.5711529 | 3 | 0.6159766 |
| 1,1,1-Trichloroethane | 0.025 | | 0.05 | | 0.1 | 3.425432 | 0.2 | 3.411367 | 0.5 | 3.411368 | 3 | 3.609574 |
| 1,1,2,2-Tetrachloroethane | 0.025 | | 0.05 | | 0.1 | 0.7300414 | 0.2 | 0.7318402 | 0.5 | 0.7544617 | 3 | 0.8690051 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 0.025 | | 0.05 | | 0.1 | 3.080905 | 0.2 | 3.021607 | 0.5 | 2.825011 | 3 | 3.008436 |
| 1,1,2-Trichloroethane | 0.025 | | 0.05 | | 0.1 | 0.3566161 | 0.2 | 0.349048 | 0.5 | 0.3485183 | 3 | 0.3808134 |
| 1,1-Dichloroethane | 0.025 | | 0.05 | | 0.1 | 2.123927 | 0.2 | 2.145415 | 0.5 | 2.038435 | 3 | 2.15819 |
| 1,1-Dichloroethylene | 0.025 | 1.68956 | 0.05 | 1.949682 | 0.1 | 1.877836 | 0.2 | 1.809408 | 0.5 | 1.774184 | 3 | 1.851395 |
| 1,2,4-Trichlorobenzene | 0.025 | | 0.05 | | 0.1 | | 0.2 | 4.587566E-02 | 0.5 | 8.934861E-02 | 3 | 0.2059795 |
| 1,2,4-Trimethylbenzene | 0.025 | | 0.05 | | 0.1 | 1.12508 | 0.2 | 1.260815 | 0.5 | 1.444108 | 3 | 1.730441 |
| 1,2-Dibromoethane | 0.025 | | 0.05 | | 0.1 | 0.5520995 | 0.2 | 0.5474189 | 0.5 | 0.5562918 | 3 | 0.6228938 |
| 1,2-Dichlorobenzene | 0.025 | | 0.05 | | 0.1 | 0.5467859 | 0.2 | 0.6057459 | 0.5 | 0.7379992 | 3 | 1.020486 |
| 1,2-Dichloroethane | 0.025 | | 0.05 | | 0.1 | 2.027475 | 0.2 | 1.97455 | 0.5 | 1.95929 | 3 | 2.021334 |
| 1,2-Dichloropropane | 0.025 | | 0.05 | | 0.1 | 0.2425652 | 0.2 | 0.23658 | 0.5 | 0.2333119 | 3 | 0.2377875 |
| 1,2-Dichlorotetrafluoroethane | 0.025 | | 0.05 | | 0.1 | 2.703432 | 0.2 | 2.711149 | 0.5 | 2.734652 | 3 | 2.868186 |
| 1,3,5-Trimethylbenzene | 0.025 | | 0.05 | | 0.1 | 1.378163 | 0.2 | 1.431125 | 0.5 | 1.536423 | 3 | 1.761684 |
| 1,3-Butadiene | 0.025 | | 0.05 | | 0.1 | 0.6378528 | 0.2 | 0.5007399 | 0.5 | 0.503555 | 3 | 0.5205629 |
| 1,3-Dichlorobenzene | 0.025 | | 0.05 | | 0.1 | 0.4679316 | 0.2 | 0.4986744 | 0.5 | 0.6136063 | 3 | 0.9167642 |
| 1,3-Dichloropropane | 0.025 | | 0.05 | | 0.1 | 0.4780494 | 0.2 | 0.458046 | 0.5 | 0.451178 | 3 | 0.5048935 |
| 1,4-Dichlorobenzene | 0.025 | | 0.05 | | 0.1 | 0.3212931 | 0.2 | 0.3820276 | 0.5 | 0.4678002 | 3 | 0.7902146 |
| 1,4-Dioxane | 0.025 | | 0.05 | | 0.1 | 0.1948808 | 0.2 | 0.1933201 | 0.5 | 0.1869593 | 3 | 0.1931201 |
| 2-Butanone | 0.025 | | 0.05 | | 0.1 | 1.80242 | 0.2 | 1.687187 | 0.5 | 1.689676 | 3 | 1.751541 |
| 2-Hexanone | 0.025 | | 0.05 | | 0.1 | 0.2555221 | 0.2 | 0.281305 | 0.5 | 0.3053715 | 3 | 0.3554359 |
| 3-Chloropropene | 0.025 | | 0.05 | | 0.1 | 0.9799991 | 0.2 | 0.9587636 | 0.5 | 0.9178889 | 3 | 0.9939028 |
| 4-Methyl-2-pentanone | 0.025 | | 0.05 | | 0.1 | 0.358424 | 0.2 | 0.3446912 | 0.5 | 0.359746 | 3 | 0.4131591 |
| Acetone | 0.025 | | 0.05 | | 0.1 | 1.969524 | 0.2 | 1.87522 | 0.5 | 1.74095 | 3 | 1.471826 |
| Acrolein | 0.025 | | 0.05 | | 0.1 | 0.3897769 | 0.2 | 0.3533793 | 0.5 | 0.3748426 | 3 | 0.3913719 |
| Acrylonitrile | 0.025 | | 0.05 | | 0.1 | 0.7223971 | 0.2 | 0.6266402 | 0.5 | 0.6583513 | 3 | 0.680862 |
| Benzene | 0.025 | | 0.05 | | 0.1 | 4.427659 | 0.2 | 4.203149 | 0.5 | 4.077926 | 3 | 4.280527 |
| Benzyl chloride | 0.025 | | 0.05 | | 0.1 | | 0.2 | 0.223463 | 0.5 | 0.3109399 | 3 | 0.7215757 |
| Bromodichloromethane | 0.025 | | 0.05 | | 0.1 | 0.6112342 | 0.2 | 0.5771457 | 0.5 | 0.5940702 | 3 | 0.6458868 |
| Bromoform | 0.025 | | 0.05 | | 0.1 | 0.5140861 | 0.2 | 0.541486 | 0.5 | 0.5916167 | 3 | 0.7717672 |

FORM VI

INITIAL CALIBRATION DATA

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.

SDG: 21A0950

Client: Advanced Cleanup Technologies, Inc.

Project: 9628-BXNY

Calibration: YA10030

Instrument: TO15 AIR2

Calibration Date: 01/26/21 07:48

| Compound | Level 01 | | Level 02 | | Level 03 | | Level 04 | | Level 05 | | Level 06 | |
|--------------------------------|----------|----------|----------|----------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| Bromomethane | 0.025 | | 0.05 | | 0.1 | 1.410659 | 0.2 | 1.256755 | 0.5 | 1.123878 | 3 | 1.167802 |
| Carbon disulfide | 0.025 | | 0.05 | | 0.1 | 3.259519 | 0.2 | 3.239685 | 0.5 | 3.170763 | 3 | 3.500906 |
| Carbon tetrachloride | 0.025 | | 0.05 | 3.476406 | 0.1 | 3.558004 | 0.2 | 3.436302 | 0.5 | 3.497501 | 3 | 3.818056 |
| Chlorobenzene | 0.025 | | 0.05 | | 0.1 | 1.06513 | 0.2 | 1.027989 | 0.5 | 1.009538 | 3 | 1.07065 |
| Chloroethane | 0.025 | | 0.05 | | 0.1 | 0.5183794 | 0.2 | 0.483776 | 0.5 | 0.4912851 | 3 | 0.4869944 |
| Chloroform | 0.025 | | 0.05 | | 0.1 | 3.176563 | 0.2 | 3.141376 | 0.5 | 3.057977 | 3 | 3.225084 |
| Chloromethane | 0.025 | | 0.05 | | 0.1 | 0.6731788 | 0.2 | 0.5730917 | 0.5 | 0.4268476 | 3 | 0.4988914 |
| cis-1,2-Dichloroethylene | 0.025 | 1.639302 | 0.05 | 1.646599 | 0.1 | 1.735341 | 0.2 | 1.609317 | 0.5 | 1.564295 | 3 | 1.609182 |
| cis-1,3-Dichloropropylene | 0.025 | | 0.05 | | 0.1 | 0.3929255 | 0.2 | 0.4017156 | 0.5 | 0.4086907 | 3 | 0.4683334 |
| Cyclohexane | 0.025 | | 0.05 | | 0.1 | 1.457893 | 0.2 | 1.444583 | 0.5 | 1.475399 | 3 | 1.528986 |
| Dibromochloromethane | 0.025 | | 0.05 | | 0.1 | 0.6215546 | 0.2 | 0.619673 | 0.5 | 0.6440579 | 3 | 0.7527804 |
| Dichlorodifluoromethane | 0.025 | | 0.05 | | 0.1 | 4.006922 | 0.2 | 3.837507 | 0.5 | 3.677406 | 3 | 3.863792 |
| Ethanol | 0.025 | | 0.05 | | 0.1 | | 0.2 | | 0.5 | | 3 | |
| Ethyl acetate | 0.025 | | 0.05 | | 0.1 | 1.875057 | 0.2 | 1.927338 | 0.5 | 1.836103 | 3 | 1.990347 |
| Ethyl Benzene | 0.025 | | 0.05 | | 0.1 | 1.469705 | 0.2 | 1.493129 | 0.5 | 1.503685 | 3 | 1.656591 |
| Hexachlorobutadiene | 0.025 | | 0.05 | | 0.1 | 0.7671693 | 0.2 | 0.7998451 | 0.5 | 0.8626803 | 3 | 0.9640472 |
| Isopropanol | 0.025 | | 0.05 | | 0.1 | 2.060419 | 0.2 | 1.815335 | 0.5 | 1.716735 | 3 | 1.642508 |
| Isopropylbenzene | 0.025 | | 0.05 | | 0.1 | 1.639762 | 0.2 | 1.710464 | 0.5 | 1.816235 | 3 | 2.068458 |
| Methyl Methacrylate | 0.025 | | 0.05 | | 0.1 | 0.2372921 | 0.2 | 0.232493 | 0.5 | 0.2354746 | 3 | 0.2772458 |
| Methyl tert-butyl ether (MTBE) | 0.025 | | 0.05 | | 0.1 | 3.677477 | 0.2 | 3.634145 | 0.5 | 3.532442 | 3 | 3.816353 |
| Methylene chloride | 0.025 | | 0.05 | | 0.1 | 2.041367 | 0.2 | 1.516118 | 0.5 | 1.139073 | 3 | 1.041112 |
| Naphthalene | 0.025 | | 0.05 | | 0.1 | 0.4697198 | 0.2 | 0.4090934 | 0.5 | 0.4624087 | 3 | 0.8830937 |
| n-Butylbenzene | 0.025 | | 0.05 | | 0.1 | 0.5775272 | 0.2 | 0.7249461 | 0.5 | 1.036054 | 3 | 1.668158 |
| n-Heptane | 0.025 | | 0.05 | | 0.1 | 1.33445 | 0.2 | 1.423736 | 0.5 | 1.385772 | 3 | 1.535934 |
| n-Hexane | 0.025 | | 0.05 | | 0.1 | 2.494255 | 0.2 | 1.856212 | 0.5 | 1.564945 | 3 | 1.486143 |
| n-Propylbenzene | 0.025 | | 0.05 | | 0.1 | 1.590797 | 0.2 | 1.702634 | 0.5 | 1.86525 | 3 | 2.239676 |
| o-Xylene | 0.025 | | 0.05 | | 0.1 | 1.139641 | 0.2 | 1.115911 | 0.5 | 1.181538 | 3 | 1.333251 |
| p- & m- Xylenes | 0.05 | | 0.1 | | 0.2 | 1.161484 | 0.4 | 1.140338 | 1 | 1.19004 | 6 | 1.298488 |
| p-Ethyltoluene | 0.025 | | 0.05 | | 0.1 | 1.165358 | 0.2 | 1.348438 | 0.5 | 1.532066 | 3 | 1.953768 |
| p-Isopropyltoluene | 0.025 | | 0.05 | | 0.1 | 1.431556 | 0.2 | 1.588328 | 0.5 | 1.83419 | 3 | 2.239238 |
| Propylene | 0.025 | | 0.05 | | 0.1 | 0.3731062 | 0.2 | 0.4034533 | 0.5 | 0.3687482 | 3 | 0.3748513 |

FORM VI

INITIAL CALIBRATION DATA

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YA10030Instrument: TO15 AIR2Calibration Date: 01/26/21 07:48

| Compound | Level 01 | | Level 02 | | Level 03 | | Level 04 | | Level 05 | | Level 06 | |
|--------------------------------------|----------|----------------------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| sec-Butylbenzene | 0.025 | | 0.05 | | 0.1 | 1.902468 | 0.2 | 1.99823 | 0.5 | 2.240514 | 3 | 2.639129 |
| Styrene | 0.025 | | 0.05 | | 0.1 | 0.669751 | 0.2 | 0.6997102 | 0.5 | 0.8053752 | 3 | 1.020018 |
| SURR: p-Bromofluorobenzene | 10 | 0.5917769 | 10 | 0.6198109 | 10 | 0.6425812 | 10 | 0.6638616 | 10 | 0.6777516 | 10 | 0.7783833 |
| tert-Butylbenzene | 0.025 | | 0.05 | | 0.1 | 1.500702 | 0.2 | 1.489469 | 0.5 | 1.649558 | 3 | 2.002725 |
| Tetrachloroethylene | 0.025 | | 0.05 | | 0.1 | 0.6638151 | 0.2 | 0.6452357 | 0.5 | 0.6234892 | 3 | 0.6486787 |
| Tetrahydrofuran | 0.025 | | 0.05 | | 0.1 | 0.9224455 | 0.2 | 0.8300019 | 0.5 | 0.8610897 | 3 | 0.92248 |
| Toluene | 0.025 | | 0.05 | | 0.1 | 1.084839 | 0.2 | 1.041284 | 0.5 | 1.047442 | 3 | 1.108024 |
| trans-1,2-Dichloroethylene | 0.025 | | 0.05 | | 0.1 | 1.696045 | 0.2 | 1.683917 | 0.5 | 1.578272 | 3 | 1.658149 |
| trans-1,3-Dichloropropylene | 0.025 | | 0.05 | | 0.1 | 0.3699496 | 0.2 | 0.3611546 | 0.5 | 0.3880453 | 3 | 0.4613246 |
| Trichloroethylene | 0.025 | 0.4670628 | 0.05 | 0.4286861 | 0.1 | 0.4038485 | 0.2 | 0.3997878 | 0.5 | 0.3831523 | 3 | 0.4075425 |
| Trichlorofluoromethane (Freon 11) | 0.025 | | 0.05 | | 0.1 | 4.200223 | 0.2 | 4.081745 | 0.5 | 3.862837 | 3 | 4.114374 |
| Vinyl acetate | 0.025 | | 0.05 | | 0.1 | 2.12869 | 0.2 | 2.069793 | 0.5 | 2.101735 | 3 | 2.381841 |
| Vinyl bromide | 0.025 | | 0.05 | | 0.1 | 1.280072 | 0.2 | 1.275967 | 0.5 | 1.212855 | 3 | 1.305876 |
| Vinyl Chloride | 0.025 | 0.9918589 | 0.05 | 0.9575005 | 0.1 | 0.7426401 | 0.2 | 0.7770665 | 0.5 | 0.7553732 | 3 | 0.7589363 |

FORM VI

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YA10030Instrument: TO15 AIR2Calibration Date: 01/26/21 07:48

| Compound | Level 07 | | Level 08 | | Level 09 | | Level 10 | | Level 11 | | Level 12 | |
|---|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|----|----------|----|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| 1,1,1,2-Tetrachloroethane | 10 | 0.581416 | 20 | 0.5856221 | 30 | 0.5335512 | 50 | 0.6696643 | | | | |
| 1,1,1-Trichloroethane | 10 | 3.378862 | 20 | 3.213151 | 30 | 2.921064 | 50 | 3.163491 | | | | |
| 1,1,2,2-Tetrachloroethane | 10 | 0.8528597 | 20 | 0.8527992 | 30 | 0.7946369 | 50 | 1.015743 | | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 10 | 2.815042 | 20 | 2.757495 | 30 | 2.515044 | 50 | 2.728945 | | | | |
| 1,1,2-Trichloroethane | 10 | 0.3666615 | 20 | 0.3494872 | 30 | 0.3129771 | 50 | 0.3600748 | | | | |
| 1,1-Dichloroethane | 10 | 2.061256 | 20 | 2.014063 | 30 | 1.854766 | 50 | 2.0735 | | | | |
| 1,1-Dichloroethylene | 10 | 1.790041 | 20 | 1.783326 | 30 | 1.645609 | 50 | 1.859642 | | | | |
| 1,2,4-Trichlorobenzene | 10 | 0.5144995 | 20 | 0.691163 | 30 | 0.7288036 | 50 | 0.9368631 | | | | |
| 1,2,4-Trimethylbenzene | 10 | 1.700383 | 20 | 1.667875 | 30 | 1.529488 | 50 | 1.884512 | | | | |
| 1,2-Dibromoethane | 10 | 0.6319934 | 20 | 0.6180941 | 30 | 0.5595337 | 50 | 0.6439212 | | | | |
| 1,2-Dichlorobenzene | 10 | 1.0758 | 20 | 1.069244 | 30 | 0.9820023 | 50 | 1.207579 | | | | |
| 1,2-Dichloroethane | 10 | 1.946935 | 20 | 1.875178 | 30 | 1.708617 | 50 | 1.906216 | | | | |
| 1,2-Dichloropropane | 10 | 0.2305082 | 20 | 0.2329614 | 30 | 0.2130162 | 50 | 0.2520294 | | | | |
| 1,2-Dichlorotetrafluoroethane | 10 | 2.644649 | 20 | 2.454798 | 30 | 2.386505 | 50 | 2.48684 | | | | |
| 1,3,5-Trimethylbenzene | 10 | 1.658946 | 20 | 1.608872 | 30 | 1.474374 | 50 | 1.796116 | | | | |
| 1,3-Butadiene | 10 | 0.5233552 | 20 | 0.4992007 | 30 | 0.4734438 | 50 | 0.5454222 | | | | |
| 1,3-Dichlorobenzene | 10 | 1.048843 | 20 | 1.078101 | 30 | 0.9979388 | 50 | 1.246324 | | | | |
| 1,3-Dichloropropane | 10 | 0.5064076 | 20 | 0.5030013 | 30 | 0.459745 | 50 | 0.5488471 | | | | |
| 1,4-Dichlorobenzene | 10 | 1.01241 | 20 | 1.067838 | 30 | 0.9998353 | 50 | 1.250273 | | | | |
| 1,4-Dioxane | 10 | 0.1944579 | 20 | 0.1995119 | 30 | 0.1842576 | 50 | 0.2180579 | | | | |
| 2-Butanone | 10 | 1.716279 | 20 | 1.743002 | 30 | 1.634339 | 50 | 1.864861 | | | | |
| 2-Hexanone | 10 | 0.3990241 | 20 | 0.3981502 | 30 | 0.3621916 | 50 | 0.4278325 | | | | |
| 3-Chloropropene | 10 | 0.9852444 | 20 | 1.005407 | 30 | 0.9502787 | 50 | 1.09447 | | | | |
| 4-Methyl-2-pentanone | 10 | 0.4159703 | 20 | 0.4308976 | 30 | 0.4022537 | 50 | 0.4904324 | | | | |
| Acetone | 10 | 1.434316 | 20 | 1.424185 | 30 | 1.317004 | 50 | 1.495717 | | | | |
| Acrolein | 10 | 0.3923452 | 20 | 0.4141052 | 30 | 0.3967547 | 50 | 0.4599698 | | | | |
| Acrylonitrile | 10 | 0.6838996 | 20 | 0.6999701 | 30 | 0.6586198 | 50 | 0.7735947 | | | | |
| Benzene | 10 | 4.008125 | 20 | 3.790132 | 30 | 3.40501 | 50 | 3.642948 | | | | |
| Benzyl chloride | 10 | 1.169678 | 20 | 1.339341 | 30 | 1.313488 | 50 | 1.681651 | | | | |
| Bromodichloromethane | 10 | 0.6406829 | 20 | 0.6337108 | 30 | 0.5783242 | 50 | 0.6793282 | | | | |
| Bromoform | 10 | 0.783714 | 20 | 0.7755094 | 30 | 0.7021665 | 50 | 0.8559322 | | | | |

FORM VI

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YA10030Instrument: TO15_AIR2Calibration Date: 01/26/21 07:48

| Compound | Level 07 | | Level 08 | | Level 09 | | Level 10 | | Level 11 | | Level 12 | |
|--------------------------------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|----|----------|----|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| Bromomethane | 10 | 1.131547 | 20 | 1.127421 | 30 | 1.050596 | 50 | 1.212408 | | | | |
| Carbon disulfide | 10 | 3.436351 | 20 | 3.392799 | 30 | 3.129129 | 50 | 3.495396 | | | | |
| Carbon tetrachloride | 10 | 3.619325 | 20 | 3.458393 | 30 | 3.115196 | 50 | 3.343215 | | | | |
| Chlorobenzene | 10 | 1.028076 | 20 | 1.0448 | 30 | 0.9613176 | 50 | 1.211787 | | | | |
| Chloroethane | 10 | 0.4776877 | 20 | 0.4953203 | 30 | 0.471101 | 50 | 0.565263 | | | | |
| Chloroform | 10 | 3.044963 | 20 | 2.952718 | 30 | 2.686411 | 50 | 2.932302 | | | | |
| Chloromethane | 10 | 0.509697 | 20 | 0.4844918 | 30 | 0.4579101 | 50 | 0.5342357 | | | | |
| cis-1,2-Dichloroethylene | 10 | 1.550769 | 20 | 1.543264 | 30 | 1.422704 | 50 | 1.619167 | | | | |
| cis-1,3-Dichloropropylene | 10 | 0.4765409 | 20 | 0.4847052 | 30 | 0.4490265 | 50 | 0.5417608 | | | | |
| Cyclohexane | 10 | 1.490509 | 20 | 1.51525 | 30 | 1.426036 | 50 | 1.627744 | | | | |
| Dibromochloromethane | 10 | 0.7637413 | 20 | 0.7423316 | 30 | 0.6669046 | 50 | 0.7599561 | | | | |
| Dichlorodifluoromethane | 10 | 3.761168 | 20 | 3.365812 | 30 | 3.033095 | 50 | 3.202049 | | | | |
| Ethanol | 10 | | 20 | | 30 | | 50 | | | | | |
| Ethyl acetate | 10 | 1.929362 | 20 | 1.97392 | 30 | 1.843368 | 50 | 2.099569 | | | | |
| Ethyl Benzene | 10 | 1.602992 | 20 | 1.619929 | 30 | 1.473947 | 50 | 1.806922 | | | | |
| Hexachlorobutadiene | 10 | 0.8355291 | 20 | 0.7289623 | 30 | 0.6130552 | 50 | 0.636539 | | | | |
| Isopropanol | 10 | 1.611729 | 20 | 1.664166 | 30 | 1.593262 | 50 | 1.788076 | | | | |
| Isopropylbenzene | 10 | 1.955893 | 20 | 1.889129 | 30 | 1.708118 | 50 | 2.03587 | | | | |
| Methyl Methacrylate | 10 | 0.2778188 | 20 | 0.2788584 | 30 | 0.2549856 | 50 | 0.300533 | | | | |
| Methyl tert-butyl ether (MTBE) | 10 | 3.694842 | 20 | 3.655046 | 30 | 3.390228 | 50 | 3.793513 | | | | |
| Methylene chloride | 10 | 0.9746953 | 20 | 0.9781346 | 30 | 0.9135102 | 50 | 1.045886 | | | | |
| Naphthalene | 10 | 1.791586 | 20 | 2.11018 | 30 | 2.035688 | 50 | 2.349137 | | | | |
| n-Butylbenzene | 10 | 1.768842 | 20 | 1.74019 | 30 | 1.574801 | 50 | 1.8739 | | | | |
| n-Heptane | 10 | 1.459586 | 20 | 1.417949 | 30 | 1.292768 | 50 | 1.425655 | | | | |
| n-Hexane | 10 | 1.407688 | 20 | 1.436321 | 30 | 1.351676 | 50 | 1.524392 | | | | |
| n-Propylbenzene | 10 | 2.191714 | 20 | 2.155994 | 30 | 1.9272 | 50 | 2.294275 | | | | |
| o-Xylene | 10 | 1.266019 | 20 | 1.253488 | 30 | 1.147425 | 50 | 1.436615 | | | | |
| p- & m- Xylenes | 20 | 1.225773 | 40 | 1.191005 | 60 | 1.046835 | 100 | 1.204829 | | | | |
| p-Ethyltoluene | 10 | 1.957115 | 20 | 1.928308 | 30 | 1.750363 | 50 | 2.092686 | | | | |
| p-Isopropyltoluene | 10 | 2.128936 | 20 | 2.053732 | 30 | 1.830759 | 50 | 2.145489 | | | | |
| Propylene | 10 | 0.3980964 | 20 | 0.3810167 | 30 | 0.3578397 | 50 | 0.403325 | | | | |

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YA10030Instrument: TO15_AIR2Calibration Date: 01/26/21 07:48

| Compound | Level 07 | | Level 08 | | Level 09 | | Level 10 | | Level 11 | | Level 12 | |
|--------------------------------------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|----|----------|----|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| sec-Butylbenzene | 10 | 2.435504 | 20 | 2.336617 | 30 | 2.07754 | 50 | 2.422593 | | | | |
| Styrene | 10 | 1.012569 | 20 | 1.001453 | 30 | 0.9127588 | 50 | 1.145732 | | | | |
| SURR: p-Bromofluorobenzene | 10 | 0.8441321 | 10 | 0.859791 | 10 | 0.8895966 | 10 | 0.9002172 | | | | |
| tert-Butylbenzene | 10 | 1.876768 | 20 | 1.825633 | 30 | 1.664738 | 50 | 2.044541 | | | | |
| Tetrachloroethylene | 10 | 0.6266036 | 20 | 0.5816288 | 30 | 0.5087127 | 50 | 0.5592684 | | | | |
| Tetrahydrofuran | 10 | 0.8932596 | 20 | 0.9230246 | 30 | 0.8710695 | 50 | 1.007943 | | | | |
| Toluene | 10 | 1.087116 | 20 | 1.073499 | 30 | 0.9724611 | 50 | 1.111755 | | | | |
| trans-1,2-Dichloroethylene | 10 | 1.598888 | 20 | 1.60735 | 30 | 1.491012 | 50 | 1.695252 | | | | |
| trans-1,3-Dichloropropylene | 10 | 0.4747695 | 20 | 0.4745293 | 30 | 0.4407643 | 50 | 0.5287012 | | | | |
| Trichloroethylene | 10 | 0.397174 | 20 | 0.3983018 | 30 | 0.3659061 | 50 | 0.436585 | | | | |
| Trichlorofluoromethane (Freon 11) | 10 | 3.844997 | 20 | 3.623695 | 30 | 3.291421 | 50 | 3.529087 | | | | |
| Vinyl acetate | 10 | 2.353557 | 20 | 2.375869 | 30 | 2.220451 | 50 | 2.500611 | | | | |
| Vinyl bromide | 10 | 1.262316 | 20 | 1.260892 | 30 | 1.177163 | 50 | 1.30907 | | | | |
| Vinyl Chloride | 10 | 0.7625025 | 20 | 0.7263405 | 30 | 0.6873932 | 50 | 0.7910524 | | | | |

FORM VI

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YA10030Instrument: TO15_AIR2Calibration Date: 01/26/21 07:48

| Compound | Mean RF | RF RSD | Mean RT | RT RSD | Linear r | Quad COD | LIMIT | Q |
|--|-----------|----------|----------|--------------|----------|----------|-------|---|
| 1,1,1,2-Tetrachloroethane | 0.5887094 | 6.742072 | 19.01263 | 0.0114587 | | | 30 | |
| 1,1,1-Trichloroethane | 3.316789 | 6.345866 | 12.57525 | 7.912986E-02 | | | 30 | |
| 1,1,2,2-Tetrachloroethane | 0.8251734 | 11.5388 | 20.8245 | 0.0100768 | | | 30 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 2.844061 | 6.575865 | 8.331125 | 0.2331768 | | | 30 | |
| 1,1,2-Trichloroethane | 0.3530245 | 5.53558 | 16.85575 | 3.204143E-02 | | | 30 | |
| 1,1-Dichloroethane | 2.058694 | 4.712681 | 10.5685 | 0.1102975 | | | 30 | |
| 1,1-Dichloroethylene | 1.803068 | 4.951396 | 8.5566 | 0.2223501 | | | 30 | |
| 1,2,4-Trichlorobenzene | 0.4589333 | 75.97585 | 27.02386 | 2.387444E-03 | | | | |
| 1,2,4-Trimethylbenzene | 1.542838 | 16.54976 | 22.55475 | 2.475774E-02 | | | 30 | |
| 1,2-Dibromoethane | 0.5915308 | 6.951902 | 18.16275 | 3.389631E-02 | | | 30 | |
| 1,2-Dichlorobenzene | 0.9057053 | 26.80984 | 24.28188 | 5.106048E-03 | | | 30 | |
| 1,2-Dichloroethane | 1.927449 | 5.316404 | 13.11138 | 5.986794E-02 | | | 30 | |
| 1,2-Dichloropropane | 0.234845 | 4.741041 | 14.50087 | 4.866836E-02 | | | 30 | |
| 1,2-Dichlorotetrafluoroethane | 2.623776 | 6.279934 | 5.340625 | 0.500558 | | | 30 | |
| 1,3,5-Trimethylbenzene | 1.580713 | 9.645799 | 21.72275 | 1.746863E-02 | | | 30 | |
| 1,3-Butadiene | 0.5255166 | 9.526717 | 5.90475 | 0.4518011 | | | 30 | |
| 1,3-Dichlorobenzene | 0.8585229 | 34.09283 | 23.38837 | 1.731669E-02 | | | | |
| 1,3-Dichloropropane | 0.488771 | 6.784468 | 17.2825 | 1.875635E-02 | | | 30 | |
| 1,4-Dichlorobenzene | 0.7864615 | 44.88766 | 23.5785 | 1.775437E-02 | | | | |
| 1,4-Dioxane | 0.1955707 | 5.243547 | 14.823 | 3.216074E-02 | | | 30 | |
| 2-Butanone | 1.736163 | 4.162562 | 11.1825 | 8.345571E-02 | | | 30 | |
| 2-Hexanone | 0.3481041 | 17.69719 | 16.86312 | 2.105124E-02 | | | 30 | |
| 3-Chloropropene | 0.9857443 | 5.271001 | 9.155125 | 0.1877687 | | | 30 | |
| 4-Methyl-2-pentanone | 0.4019468 | 11.87344 | 15.40063 | 2.243193E-02 | | | 30 | |
| Acetone | 1.591093 | 14.9728 | 8.19325 | 0.2336922 | | | 30 | |
| Acrolein | 0.3965682 | 7.83366 | 8.08125 | 0.2306581 | | | 30 | |
| Acrylonitrile | 0.6880419 | 6.552496 | 9.392 | 0.1400769 | | | 30 | |
| Benzene | 3.979434 | 8.659068 | 13.23963 | 5.189978E-02 | | | 30 | |
| Benzyl chloride | 0.9657338 | 57.55983 | 23.72957 | 1.971246E-02 | | | | |
| Bromodichloromethane | 0.6200479 | 5.823982 | 14.86175 | 4.033456E-02 | | | 30 | |
| Bromoform | 0.6920348 | 18.36848 | 20.57175 | 2.173355E-02 | | | 30 | |
| Bromomethane | 1.185133 | 9.303023 | 6.70075 | 0.4422791 | | | 30 | |

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YA10030Instrument: TO15_AIR2Calibration Date: 01/26/21 07:48

| Compound | Mean RF | RF RSD | Mean RT | RT RSD | Linear r | Quad COD | LIMIT | Q |
|--------------------------------|-----------|-----------------|----------|--------------|----------|-----------|-------|---|
| Carbon disulfide | 3.328068 | 4.407572 | 9.435125 | 0.172099 | | | 30 | |
| Carbon tetrachloride | 3.480266 | 5.51077 | 13.03944 | 7.550187E-02 | | | 30 | |
| Chlorobenzene | 1.052411 | 6.934101 | 18.96012 | 1.199453E-02 | | | 30 | |
| Chloroethane | 0.4987259 | 6.085233 | 6.879 | 0.3616734 | | | 30 | |
| Chloroform | 3.027174 | 5.6732 | 11.83712 | 9.981308E-02 | | | 30 | |
| Chloromethane | 0.519793 | 14.69297 | 5.554875 | 0.482107 | | | 30 | |
| cis-1,2-Dichloroethylene | 1.593994 | 5.146297 | 11.5603 | 9.577007E-02 | | | 30 | |
| cis-1,3-Dichloropropylene | 0.4529623 | 11.16112 | 15.7495 | 3.843104E-02 | | | 30 | |
| Cyclohexane | 1.4958 | 4.247787 | 12.74425 | 5.676724E-02 | | | 30 | |
| Dibromochloromethane | 0.696375 | 9.236684 | 17.79038 | 5.579376E-03 | | | 30 | |
| Dichlorodifluoromethane | 3.593469 | 9.741577 | 5.071625 | 0.6085702 | | | 30 | |
| Ethanol | | | | | | | 30 | |
| Ethyl acetate | 1.934383 | 4.520225 | 11.48312 | 8.049565E-02 | | | 30 | |
| Ethyl Benzene | 1.578363 | 7.429017 | 19.06987 | 8.385096E-03 | | | 30 | |
| Hexachlorobutadiene | 0.7759784 | 15.03483 | 27.29825 | 1.757873E-02 | | | 30 | |
| Isopropanol | 1.736529 | 8.823526 | 7.982625 | 0.211466 | | 0.9991326 | 0.99 | |
| Isopropylbenzene | 1.852991 | 8.651179 | 20.64937 | 1.930163E-02 | | | 30 | |
| Methyl Methacrylate | 0.2618377 | 9.667602 | 14.51288 | 3.784646E-02 | | | 30 | |
| Methyl tert-butyl ether (MTBE) | 3.649256 | 3.771654 | 9.717125 | 0.1303953 | | | 30 | |
| Methylene chloride | 1.086933 | 18.59915 | 9.301857 | 0.1767361 | | | 30 | |
| Naphthalene | 1.313863 | 63.65916 | 27.41525 | 0.0130803 | | | | |
| n-Butylbenzene | 1.370552 | <u>37.37237</u> | 24.021 | 1.602341E-02 | | | | |
| n-Heptane | 1.409481 | 5.28559 | 13.25725 | 6.229228E-02 | | | 30 | |
| n-Hexane | 1.640204 | 23.01786 | 10.1955 | 0.1550243 | | | 30 | |
| n-Propylbenzene | 1.995942 | 13.17873 | 21.41425 | 1.388813E-02 | | | 30 | |
| o-Xylene | 1.234236 | 8.949367 | 19.99587 | 1.820239E-02 | | | 30 | |
| p- & m- Xylenes | 1.182349 | 6.113822 | 19.21362 | 2.252799E-02 | | | 30 | |
| p-Ethyltoluene | 1.716013 | 19.38963 | 21.63662 | 1.104796E-02 | | | 30 | |
| p-Isopropyltoluene | 1.906529 | 15.0558 | 23.22113 | 1.651333E-02 | | | 30 | |
| Propylene | 0.3825546 | 4.486574 | 4.990625 | 0.5658637 | | | 30 | |
| sec-Butylbenzene | 2.256574 | 11.06959 | 22.95613 | 7.907417E-03 | | | 30 | |
| Styrene | 0.9084209 | 18.58071 | 20.01463 | 2.539488E-02 | | | 30 | |

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YA10030Instrument: TO15_AIR2Calibration Date: 01/26/21 07:48

| Compound | Mean RF | RF RSD | Mean RT | RT RSD | Linear r | Quad COD | LIMIT | Q |
|-----------------------------------|-----------|----------|----------|--------------|----------|----------|-------|---|
| SURR: p-Bromofluorobenzene | 0.7640139 | 14.86116 | 21.06 | 1.653991E-02 | | | 30 | |
| tert-Butylbenzene | 1.756767 | 12.15981 | 22.48463 | 1.531508E-02 | | | 30 | |
| Tetrachloroethylene | 0.607179 | 8.729419 | 17.59625 | 2.846063E-02 | | | 30 | |
| Tetrahydrofuran | 0.9039142 | 5.967066 | 12.169 | 0.1146554 | | | 30 | |
| Toluene | 1.065802 | 4.258662 | 16.40525 | 3.746255E-02 | | | 30 | |
| trans-1,2-Dichloroethylene | 1.626111 | 4.390033 | 9.90575 | 0.1555803 | | | 30 | |
| trans-1,3-Dichloropropylene | 0.4374048 | 13.52361 | 16.56075 | 8.278457E-03 | | | 30 | |
| Trichloroethylene | 0.4088047 | 7.0201 | 14.2715 | 4.410282E-02 | | | 30 | |
| Trichlorofluoromethane (Freon 11) | 3.818547 | 8.30935 | 7.45575 | 0.2956324 | | | 30 | |
| Vinyl acetate | 2.266568 | 6.968041 | 10.52075 | 0.1233509 | | | 30 | |
| Vinyl bromide | 1.260526 | 3.580491 | 7.33225 | 0.3302443 | | | 30 | |
| Vinyl Chloride | 0.7950664 | 12.46941 | 5.8281 | 0.4904991 | | | 30 | |

FORM VI

INITIAL CALIBRATION DATA

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.

SDG: 21A0950

Client: Advanced Cleanup Technologies, Inc.

Project: 9628-BXNY

Calibration: YB10001

Instrument: 5975C

Calibration Date: 01/28/21 00:14

| Compound | Level 01 | | Level 02 | | Level 03 | | Level 04 | | Level 05 | | Level 06 | |
|---|----------|----------|----------|----------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| 1,1,1,2-Tetrachloroethane | 0.025 | | 0.05 | | 0.1 | 0.5799111 | 0.2 | 0.4966064 | 0.5 | 0.48804 | 3 | 0.5692947 |
| 1,1,1-Trichloroethane | 0.025 | | 0.05 | | 0.1 | 1.909587 | 0.2 | 1.581737 | 0.5 | 1.501309 | 3 | 1.552766 |
| 1,1,2,2-Tetrachloroethane | 0.025 | | 0.05 | | 0.1 | 0.7644537 | 0.2 | 0.6219577 | 0.5 | 0.6096898 | 3 | 1.20386 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 0.025 | | 0.05 | | 0.1 | 2.08478 | 0.2 | 1.843482 | 0.5 | 1.732745 | 3 | 1.705828 |
| 1,1,2-Trichloroethane | 0.025 | | 0.05 | | 0.1 | 0.3562615 | 0.2 | 0.3067934 | 0.5 | 0.2919996 | 3 | 0.3922658 |
| 1,1-Dichloroethane | 0.025 | | 0.05 | | 0.1 | 1.72491 | 0.2 | 1.579061 | 0.5 | 1.576933 | 3 | 1.669226 |
| 1,1-Dichloroethylene | 0.025 | 1.437436 | 0.05 | 1.247322 | 0.1 | 1.483575 | 0.2 | 1.385744 | 0.5 | 1.36061 | 3 | 1.351523 |
| 1,2,4-Trichlorobenzene | 0.025 | | 0.05 | | 0.1 | 0.1061586 | 0.2 | | 0.5 | | 3 | 0.1966694 |
| 1,2,4-Trimethylbenzene | 0.025 | | 0.05 | | 0.1 | 0.5396473 | 0.2 | 0.3423459 | 0.5 | 0.2737645 | 3 | 1.542911 |
| 1,2-Dibromoethane | 0.025 | | 0.05 | | 0.1 | 0.4788165 | 0.2 | 0.3635029 | 0.5 | 0.3614262 | 3 | 0.5653001 |
| 1,2-Dichlorobenzene | 0.025 | | 0.05 | | 0.1 | 0.5074921 | 0.2 | 0.2700729 | 0.5 | 0.2459977 | 3 | 0.911476 |
| 1,2-Dichloroethane | 0.025 | | 0.05 | | 0.1 | 1.105213 | 0.2 | 1.003193 | 0.5 | 1.01617 | 3 | 1.067533 |
| 1,2-Dichloropropane | 0.025 | | 0.05 | | 0.1 | 0.3576031 | 0.2 | 0.3192678 | 0.5 | 0.3204705 | 3 | 0.374992 |
| 1,2-Dichlorotetrafluoroethane | 0.025 | | 0.05 | | 0.1 | 2.380879 | 0.2 | 2.548322 | 0.5 | 2.654355 | 3 | 2.612255 |
| 1,3,5-Trimethylbenzene | 0.025 | | 0.05 | | 0.1 | 0.6786135 | 0.2 | 0.464303 | 0.5 | 0.3920613 | 3 | 1.540981 |
| 1,3-Butadiene | 0.025 | | 0.05 | | 0.1 | 0.8510759 | 0.2 | 0.5537627 | 0.5 | 0.5647104 | 3 | 0.5644328 |
| 1,3-Dichlorobenzene | 0.025 | | 0.05 | | 0.1 | 0.3762618 | 0.2 | 0.2222615 | 0.5 | 0.2234176 | 3 | 0.8244748 |
| 1,3-Dichloropropane | 0.025 | | 0.05 | | 0.1 | 0.4501733 | 0.2 | 0.3811288 | 0.5 | 0.3688517 | 3 | 0.5084904 |
| 1,4-Dichlorobenzene | 0.025 | | 0.05 | | 0.1 | 0.3346932 | 0.2 | 0.1810044 | 0.5 | 0.1597112 | 3 | 0.7197025 |
| 1,4-Dioxane | 0.025 | | 0.05 | | 0.1 | 0.1106819 | 0.2 | | 0.5 | | 3 | 0.2035764 |
| 2-Butanone | 0.025 | | 0.05 | | 0.1 | 0.8292656 | 0.2 | 0.7551053 | 0.5 | 0.80472 | 3 | 2.426894 |
| 2-Hexanone | 0.025 | | 0.05 | | 0.1 | | 0.2 | | 0.5 | | 3 | 0.7738453 |
| 3-Chloropropene | 0.025 | | 0.05 | | 0.1 | 1.184394 | 0.2 | 1.257476 | 0.5 | 1.220157 | 3 | 1.343808 |
| 4-Methyl-2-pentanone | 0.025 | | 0.05 | | 0.1 | | 0.2 | | 0.5 | | 3 | 0.9476634 |
| Acetone | 0.025 | | 0.05 | | 0.1 | 1.67157 | 0.2 | 1.519926 | 0.5 | 1.42402 | 3 | 1.519894 |
| Acrolein | 0.025 | | 0.05 | | 0.1 | 0.3024994 | 0.2 | 0.2861037 | 0.5 | 0.2947227 | 3 | 0.4555416 |
| Acrylonitrile | 0.025 | | 0.05 | | 0.1 | 0.7986838 | 0.2 | 0.659362 | 0.5 | 0.7195354 | 3 | 0.8872426 |
| Benzene | 0.025 | | 0.05 | | 0.1 | 3.561954 | 0.2 | 2.957062 | 0.5 | 2.680632 | 3 | 2.88207 |
| Benzyl chloride | 0.025 | | 0.05 | | 0.1 | | 0.2 | | 0.5 | | 3 | 0.2154009 |
| Bromodichloromethane | 0.025 | | 0.05 | | 0.1 | 0.5615931 | 0.2 | 0.5204759 | 0.5 | 0.519151 | 3 | 0.6031199 |
| Bromoform | 0.025 | | 0.05 | | 0.1 | 0.5713364 | 0.2 | 0.4452838 | 0.5 | 0.4452373 | 3 | 0.7097883 |

FORM VI

INITIAL CALIBRATION DATA

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.

SDG: 21A0950

Client: Advanced Cleanup Technologies, Inc.

Project: 9628-BXNY

Calibration: YB10001

Instrument: 5975C

Calibration Date: 01/28/21 00:14

| Compound | Level 01 | | Level 02 | | Level 03 | | Level 04 | | Level 05 | | Level 06 | |
|--------------------------------|----------|----------|----------|----------|----------|--------------|----------|--------------|----------|--------------|----------|-----------|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| Bromomethane | 0.025 | | 0.05 | | 0.1 | 1.078425 | 0.2 | 0.9579968 | 0.5 | 0.9288268 | 3 | 0.8357954 |
| Carbon disulfide | 0.025 | | 0.05 | | 0.1 | 3.023335 | 0.2 | 2.607458 | 0.5 | 2.570527 | 3 | 2.662203 |
| Carbon tetrachloride | 0.025 | 1.972638 | 0.05 | 1.304486 | 0.1 | 1.82685 | 0.2 | 1.554985 | 0.5 | 1.482496 | 3 | 1.510631 |
| Chlorobenzene | 0.025 | | 0.05 | | 0.1 | 1.101757 | 0.2 | 0.9051977 | 0.5 | 0.859643 | 3 | 0.9940896 |
| Chloroethane | 0.025 | | 0.05 | | 0.1 | 0.5462058 | 0.2 | 0.5057502 | 0.5 | 0.4800809 | 3 | 0.4352357 |
| Chloroform | 0.025 | | 0.05 | | 0.1 | 1.970987 | 0.2 | 1.757313 | 0.5 | 1.65817 | 3 | 1.75652 |
| Chloromethane | 0.025 | | 0.05 | | 0.1 | 0.8501276 | 0.2 | 0.7534158 | 0.5 | 0.7278626 | 3 | 0.6150147 |
| cis-1,2-Dichloroethylene | 0.025 | 1.261899 | 0.05 | 1.070051 | 0.1 | 1.32948 | 0.2 | 1.205662 | 0.5 | 1.222624 | 3 | 1.293738 |
| cis-1,3-Dichloropropylene | 0.025 | | 0.05 | | 0.1 | 0.4271649 | 0.2 | 0.3668663 | 0.5 | 0.3617917 | 3 | 0.4708267 |
| Cyclohexane | 0.025 | | 0.05 | | 0.1 | 1.792475 | 0.2 | 1.622428 | 0.5 | 1.566138 | 3 | 1.649286 |
| Dibromochloromethane | 0.025 | | 0.05 | | 0.1 | 0.5503908 | 0.2 | 0.453293 | 0.5 | 0.44498013 | 3 | 0.6147541 |
| Dichlorodifluoromethane | 0.025 | | 0.05 | | 0.1 | 2.909305 | 0.2 | 2.578594 | 0.5 | 2.316022 | 3 | 1.90122 |
| Ethanol | 0.025 | | 0.05 | | 0.1 | 0.3520467 | 0.2 | 0.5389788 | 0.5 | 0.5374464 | 3 | 0.4652192 |
| Ethyl acetate | 0.025 | | 0.05 | | 0.1 | 0.9992437 | 0.2 | 0.8253641 | 0.5 | 0.8554853 | 3 | 2.710518 |
| Ethyl Benzene | 0.025 | | 0.05 | | 0.1 | 1.440085 | 0.2 | 1.22039 | 0.5 | 1.140402 | 3 | 1.581998 |
| Hexachlorobutadiene | 0.025 | | 0.05 | | 0.1 | 0.6043304 | 0.2 | 0.1903677 | 0.5 | 0.1094953 | 3 | 0.7302993 |
| Isopropanol | 0.025 | | 0.05 | | 0.1 | 1.211657 | 0.2 | 1.037407 | 0.5 | 1.009817 | 3 | 2.140223 |
| Isopropylbenzene | 0.025 | | 0.05 | | 0.1 | 1.105671 | 0.2 | 0.8222154 | 0.5 | 0.7833892 | 3 | 1.785809 |
| Methyl Methacrylate | 0.025 | | 0.05 | | 0.1 | 8.894798E-02 | 0.2 | 7.765628E-02 | 0.5 | 8.041098E-02 | 3 | 0.3445318 |
| Methyl tert-butyl ether (MTBE) | 0.025 | | 0.05 | | 0.1 | 1.690535 | 0.2 | 1.508803 | 0.5 | 1.544241 | 3 | 2.165477 |
| Methylene chloride | 0.025 | | 0.05 | | 0.1 | 9.393839 | 0.2 | 5.248707 | 0.5 | 2.868581 | 3 | 1.555752 |
| Naphthalene | 0.025 | | 0.05 | | 0.1 | | 0.2 | | 0.5 | | 3 | 0.8465911 |
| n-Butylbenzene | 0.025 | | 0.05 | | 0.1 | | 0.2 | | 0.5 | | 3 | 1.786426 |
| n-Heptane | 0.025 | | 0.05 | | 0.1 | 2.002992 | 0.2 | 1.927961 | 0.5 | 1.935991 | 3 | 2.061876 |
| n-Hexane | 0.025 | | 0.05 | | 0.1 | 1.643359 | 0.2 | 1.465155 | 0.5 | 1.443327 | 3 | 1.49971 |
| n-Propylbenzene | 0.025 | | 0.05 | | 0.1 | 1.011443 | 0.2 | 0.733381 | 0.5 | 0.6768539 | 3 | 2.289626 |
| o-Xylene | 0.025 | | 0.05 | | 0.1 | 0.9793806 | 0.2 | 0.7941255 | 0.5 | 0.7177182 | 3 | 1.235213 |
| p- & m- Xylenes | 0.05 | | 0.1 | | 0.2 | 1.025796 | 0.4 | 0.8536117 | 1 | 0.8286805 | 6 | 1.228453 |
| p-Ethyltoluene | 0.025 | | 0.05 | | 0.1 | 0.6516777 | 0.2 | 0.4772361 | 0.5 | 0.4348378 | 3 | 1.767332 |
| p-Isopropyltoluene | 0.025 | | 0.05 | | 0.1 | 0.5345211 | 0.2 | 0.3076431 | 0.5 | 0.2452381 | 3 | 2.077021 |
| Propylene | 0.025 | | 0.05 | | 0.1 | 0.7873045 | 0.2 | 0.7199056 | 0.5 | 0.6607512 | 3 | 0.6174036 |

FORM VI

INITIAL CALIBRATION DATA

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.

SDG: 21A0950

Client: Advanced Cleanup Technologies, Inc.

Project: 9628-BXNY

Calibration: YB10001

Instrument: 5975C

Calibration Date: 01/28/21 00:14

| Compound | Level 01 | | Level 02 | | Level 03 | | Level 04 | | Level 05 | | Level 06 | |
|--------------------------------------|----------|----------------------|----------|----------------------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| sec-Butylbenzene | 0.025 | | 0.05 | | 0.1 | 0.8866433 | 0.2 | 0.546466 | 0.5 | 0.4269531 | 3 | 2.500149 |
| Styrene | 0.025 | | 0.05 | | 0.1 | 0.52576 | 0.2 | 0.3972383 | 0.5 | 0.3690359 | 3 | 0.9446457 |
| SURR: p-Bromofluorobenzene | 10 | 0.4013762 | 10 | 0.4263027 | 10 | 0.4457645 | 10 | 0.4495534 | 10 | 0.4545208 | 10 | 0.566851 |
| tert-Butylbenzene | 0.025 | | 0.05 | | 0.1 | 0.6472971 | 0.2 | 0.4442304 | 0.5 | 0.3213871 | 3 | 1.745234 |
| Tetrachloroethylene | 0.025 | | 0.05 | | 0.1 | 0.4504417 | 0.2 | 0.3735505 | 0.5 | 0.3781817 | 3 | 0.4457019 |
| Tetrahydrofuran | 0.025 | | 0.05 | | 0.1 | 0.6514643 | 0.2 | 0.5337692 | 0.5 | 0.527207 | 3 | 1.408516 |
| Toluene | 0.025 | | 0.05 | | 0.1 | 1.013175 | 0.2 | 0.896666 | 0.5 | 0.8747099 | 3 | 0.9990909 |
| trans-1,2-Dichloroethylene | 0.025 | | 0.05 | | 0.1 | 1.329006 | 0.2 | 1.248606 | 0.5 | 1.2254 | 3 | 1.312228 |
| trans-1,3-Dichloropropylene | 0.025 | | 0.05 | | 0.1 | 0.3085006 | 0.2 | 0.251276 | 0.5 | 0.2555646 | 3 | 0.4111309 |
| Trichloroethylene | 0.025 | 0.5293336 | 0.05 | 0.3681543 | 0.1 | 0.4097912 | 0.2 | 0.3719753 | 0.5 | 0.3442282 | 3 | 0.3770084 |
| Trichlorofluoromethane (Freon 11) | 0.025 | | 0.05 | | 0.1 | 2.184823 | 0.2 | 1.873331 | 0.5 | 1.822802 | 3 | 1.715396 |
| Vinyl acetate | 0.025 | | 0.05 | | 0.1 | 1.034093 | 0.2 | 0.5322205 | 0.5 | 0.3595518 | 3 | 2.886531 |
| Vinyl bromide | 0.025 | | 0.05 | | 0.1 | 0.9662912 | 0.2 | 0.8143817 | 0.5 | 0.7799233 | 3 | 0.730707 |
| Vinyl Chloride | 0.025 | 1.229984 | 0.05 | 1.017383 | 0.1 | 1.411269 | 0.2 | 0.8090314 | 0.5 | 0.7805402 | 3 | 0.7183967 |

FORM VI

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YB10001Instrument: 5975CCalibration Date: 01/28/21 00:14

| Compound | Level 07 | | Level 08 | | Level 09 | | Level 10 | | Level 11 | | Level 12 | |
|---|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|----|----------|----|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| 1,1,1,2-Tetrachloroethane | 10 | 0.5452751 | 20 | 0.551284 | 30 | 0.5552604 | 50 | 0.5149272 | | | | |
| 1,1,1-Trichloroethane | 10 | 1.565552 | 20 | 1.731954 | 30 | 1.838435 | 50 | 1.968982 | | | | |
| 1,1,2,2-Tetrachloroethane | 10 | 1.074549 | 20 | 1.046345 | 30 | 1.040907 | 50 | 0.9499997 | | | | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 10 | 1.722274 | 20 | 1.865295 | 30 | 1.922386 | 50 | 1.894504 | | | | |
| 1,1,2-Trichloroethane | 10 | 0.3963627 | 20 | 0.390164 | 30 | 0.3697174 | 50 | 0.3164285 | | | | |
| 1,1-Dichloroethane | 10 | 1.600395 | 20 | 1.651487 | 30 | 1.671907 | 50 | 1.717412 | | | | |
| 1,1-Dichloroethylene | 10 | 1.333851 | 20 | 1.384456 | 30 | 1.416557 | 50 | 1.449158 | | | | |
| 1,2,4-Trichlorobenzene | 10 | 0.4085062 | 20 | 0.5908362 | 30 | 0.6774839 | 50 | 0.7006932 | | | | |
| 1,2,4-Trimethylbenzene | 10 | 1.435829 | 20 | 1.446346 | 30 | 1.458105 | 50 | 1.375326 | | | | |
| 1,2-Dibromoethane | 10 | 0.5847049 | 20 | 0.6051196 | 30 | 0.6063129 | 50 | 0.5787043 | | | | |
| 1,2-Dichlorobenzene | 10 | 0.9436641 | 20 | 0.9661151 | 30 | 0.9594406 | 50 | 0.9008386 | | | | |
| 1,2-Dichloroethane | 10 | 1.011332 | 20 | 1.055533 | 30 | 1.111644 | 50 | 1.184971 | | | | |
| 1,2-Dichloropropane | 10 | 0.3497633 | 20 | 0.3460521 | 30 | 0.3309332 | 50 | 0.2890553 | | | | |
| 1,2-Dichlorotetrafluoroethane | 10 | 2.438996 | 20 | 2.276618 | 30 | 2.088586 | 50 | 1.885481 | | | | |
| 1,3,5-Trimethylbenzene | 10 | 1.360943 | 20 | 1.372565 | 30 | 1.366753 | 50 | 1.322004 | | | | |
| 1,3-Butadiene | 10 | 0.5347713 | 20 | 0.5196894 | 30 | 0.4863144 | 50 | 0.5135756 | | | | |
| 1,3-Dichlorobenzene | 10 | 0.9039347 | 20 | 0.9390727 | 30 | 0.9575492 | 50 | 0.9098892 | | | | |
| 1,3-Dichloropropane | 10 | 0.5051632 | 20 | 0.5047014 | 30 | 0.5061467 | 50 | 0.4993227 | | | | |
| 1,4-Dichlorobenzene | 10 | 0.8442899 | 20 | 0.9099712 | 30 | 0.9374576 | 50 | 0.9023117 | | | | |
| 1,4-Dioxane | 10 | 0.1948845 | 20 | 0.1998587 | 30 | 0.2013786 | 50 | 0.197518 | | | | |
| 2-Butanone | 10 | 2.213659 | 20 | 2.279497 | 30 | 2.268154 | 50 | 2.294415 | | | | |
| 2-Hexanone | 10 | 0.7902806 | 20 | 0.7870759 | 30 | 0.7517169 | 50 | 0.6703202 | | | | |
| 3-Chloropropene | 10 | 1.276337 | 20 | 1.29752 | 30 | 1.31881 | 50 | 1.330082 | | | | |
| 4-Methyl-2-pentanone | 10 | 0.8582218 | 20 | 0.8216563 | 30 | 0.8079306 | 50 | 0.7762045 | | | | |
| Acetone | 10 | 1.373515 | 20 | 1.416296 | 30 | 1.426911 | 50 | 1.438425 | | | | |
| Acrolein | 10 | 0.4571573 | 20 | 0.4774965 | 30 | 0.4755082 | 50 | 0.4856225 | | | | |
| Acrylonitrile | 10 | 0.8384256 | 20 | 0.8553916 | 30 | 0.8512328 | 50 | 0.8627862 | | | | |
| Benzene | 10 | 2.841214 | 20 | 2.956513 | 30 | 2.858826 | 50 | 2.563769 | | | | |
| Benzyl chloride | 10 | 0.2800357 | 20 | 0.3054501 | 30 | 0.3032601 | 50 | 0.2824589 | | | | |
| Bromodichloromethane | 10 | 0.5855758 | 20 | 0.6022534 | 30 | 0.611936 | 50 | 0.6026798 | | | | |
| Bromoform | 10 | 0.748161 | 20 | 0.7717959 | 30 | 0.7947007 | 50 | 0.743751 | | | | |

FORM VI

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YB10001Instrument: 5975CCalibration Date: 01/28/21 00:14

| Compound | Level 07 | | Level 08 | | Level 09 | | Level 10 | | Level 11 | | Level 12 | |
|--------------------------------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|----|----------|----|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| Bromomethane | 10 | 0.8348745 | 20 | 0.9242786 | 30 | 0.9614923 | 50 | 0.994344 | | | | |
| Carbon disulfide | 10 | 2.62618 | 20 | 2.760641 | 30 | 2.768882 | 50 | 2.810557 | | | | |
| Carbon tetrachloride | 10 | 1.554196 | 20 | 1.755943 | 30 | 1.919068 | 50 | 2.066129 | | | | |
| Chlorobenzene | 10 | 0.9576976 | 20 | 0.966864 | 30 | 0.9688196 | 50 | 0.9059595 | | | | |
| Chloroethane | 10 | 0.4384956 | 20 | 0.4723188 | 30 | 0.4782532 | 50 | 0.4838987 | | | | |
| Chloroform | 10 | 1.764125 | 20 | 1.873154 | 30 | 1.951718 | 50 | 2.028737 | | | | |
| Chloromethane | 10 | 0.6082463 | 20 | 0.5842623 | 30 | 0.5496103 | 50 | 0.5839539 | | | | |
| cis-1,2-Dichloroethylene | 10 | 1.221286 | 20 | 1.272999 | 30 | 1.300095 | 50 | 1.352499 | | | | |
| cis-1,3-Dichloropropylene | 10 | 0.4647274 | 20 | 0.4777595 | 30 | 0.4811049 | 50 | 0.4767408 | | | | |
| Cyclohexane | 10 | 1.573725 | 20 | 1.64601 | 30 | 1.661884 | 50 | 1.683584 | | | | |
| Dibromochloromethane | 10 | 0.6381664 | 20 | 0.6613106 | 30 | 0.6667918 | 50 | 0.6371418 | | | | |
| Dichlorodifluoromethane | 10 | 2.067146 | 20 | 2.243076 | 30 | 2.278772 | 50 | 2.286868 | | | | |
| Ethanol | 10 | 0.4360186 | 20 | 0.4476957 | 30 | 0.4475017 | 50 | 0.450724 | | | | |
| Ethyl acetate | 10 | 2.451693 | 20 | 2.534473 | 30 | 2.562783 | 50 | 2.556619 | | | | |
| Ethyl Benzene | 10 | 1.511088 | 20 | 1.506664 | 30 | 1.491921 | 50 | 1.391992 | | | | |
| Hexachlorobutadiene | 10 | 0.795959 | 20 | 0.8351738 | 30 | 0.7942857 | 50 | 0.6905302 | | | | |
| Isopropanol | 10 | 2.031593 | 20 | 2.03248 | 30 | 2.067289 | 50 | 2.070019 | | | | |
| Isopropylbenzene | 10 | 1.681823 | 20 | 1.658801 | 30 | 1.673885 | 50 | 1.548061 | | | | |
| Methyl Methacrylate | 10 | 0.330449 | 20 | 0.3412392 | 30 | 0.3322276 | 50 | 0.2946065 | | | | |
| Methyl tert-butyl ether (MTBE) | 10 | 2.042355 | 20 | 2.197956 | 30 | 2.292665 | 50 | 2.362516 | | | | |
| Methylene chloride | 10 | 1.296804 | 20 | 1.277804 | 30 | 1.259062 | 50 | 1.250825 | | | | |
| Naphthalene | 10 | 1.484326 | 20 | 1.83963 | 30 | 1.957516 | 50 | 1.900501 | | | | |
| n-Butylbenzene | 10 | 1.782706 | 20 | 1.807827 | 30 | 1.766656 | 50 | 1.584042 | | | | |
| n-Heptane | 10 | 1.99422 | 20 | 2.058956 | 30 | 1.997779 | 50 | 1.855054 | | | | |
| n-Hexane | 10 | 1.459078 | 20 | 1.505232 | 30 | 1.526851 | 50 | 1.536265 | | | | |
| n-Propylbenzene | 10 | 2.098989 | 20 | 2.115695 | 30 | 2.0732 | 50 | 1.958271 | | | | |
| o-Xylene | 10 | 1.192228 | 20 | 1.205139 | 30 | 1.177429 | 50 | 1.061583 | | | | |
| p- & m- Xylenes | 20 | 1.171805 | 40 | 1.156534 | 60 | 1.11119 | 100 | 0.9816939 | | | | |
| p-Ethyltoluene | 10 | 1.702453 | 20 | 1.712024 | 30 | 1.700738 | 50 | 1.61259 | | | | |
| p-Isopropyltoluene | 10 | 1.938093 | 20 | 1.922974 | 30 | 1.870284 | 50 | 1.722232 | | | | |
| Propylene | 10 | 0.6159713 | 20 | 0.6310291 | 30 | 0.6060071 | 50 | 0.5997247 | | | | |

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YB10001Instrument: 5975CCalibration Date: 01/28/21 00:14

| Compound | Level 07 | | Level 08 | | Level 09 | | Level 10 | | Level 11 | | Level 12 | |
|--------------------------------------|----------|-----------|----------|-----------|----------|-----------|----------|-----------|----------|----|----------|----|
| | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF | ppbv | RF |
| sec-Butylbenzene | 10 | 2.258649 | 20 | 2.252537 | 30 | 2.178913 | 50 | 1.99857 | | | | |
| Styrene | 10 | 0.9694748 | 20 | 0.9772832 | 30 | 0.9327315 | 50 | 0.8389865 | | | | |
| SURR: p-Bromofluorobenzene | 10 | 0.5839337 | 10 | 0.5798507 | 10 | 0.5866482 | 10 | 0.6030043 | | | | |
| tert-Butylbenzene | 10 | 1.626732 | 20 | 1.624493 | 30 | 1.642866 | 50 | 1.545519 | | | | |
| Tetrachloroethylene | 10 | 0.4617425 | 20 | 0.4694104 | 30 | 0.460528 | 50 | 0.430886 | | | | |
| Tetrahydrofuran | 10 | 1.271495 | 20 | 1.301304 | 30 | 1.284318 | 50 | 1.300876 | | | | |
| Toluene | 10 | 1.019481 | 20 | 1.014892 | 30 | 1.000745 | 50 | 0.9509457 | | | | |
| trans-1,2-Dichloroethylene | 10 | 1.293001 | 20 | 1.340303 | 30 | 1.385524 | 50 | 1.42023 | | | | |
| trans-1,3-Dichloropropylene | 10 | 0.4088256 | 20 | 0.4166435 | 30 | 0.4208882 | 50 | 0.4252092 | | | | |
| Trichloroethylene | 10 | 0.3658884 | 20 | 0.3809611 | 30 | 0.3862894 | 50 | 0.372664 | | | | |
| Trichlorofluoromethane (Freon 11) | 10 | 1.773389 | 20 | 1.989793 | 30 | 2.158117 | 50 | 2.231661 | | | | |
| Vinyl acetate | 10 | 2.621604 | 20 | 2.840886 | 30 | 2.969136 | 50 | 3.065108 | | | | |
| Vinyl bromide | 10 | 0.7556302 | 20 | 0.8536067 | 30 | 0.9016156 | 50 | 0.9370669 | | | | |
| Vinyl Chloride | 10 | 0.6865145 | 20 | 0.6667866 | 30 | 0.6203611 | 50 | 0.5894503 | | | | |

FORM VI

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YB10001Instrument: 5975CCalibration Date: 01/28/21 00:14

| Compound | Mean RF | RF RSD | Mean RT | RT RSD | Linear r | Quad COD | LIMIT | Q |
|--|-----------|----------|----------|--------------|----------|----------|-------|---|
| 1,1,1,2-Tetrachloroethane | 0.5375749 | 6.290038 | 19.214 | 1.677045E-02 | | | 30 | |
| 1,1,1-Trichloroethane | 1.70629 | 10.60869 | 12.80675 | 3.099859E-02 | | | 30 | |
| 1,1,2,2-Tetrachloroethane | 0.9139702 | 24.28896 | 21.02575 | 1.287143E-02 | | | 30 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 1.846412 | 6.892223 | 8.61575 | 3.125452E-02 | | | 30 | |
| 1,1,2-Trichloroethane | 0.3524991 | 11.88448 | 17.0635 | 2.318274E-02 | | | 30 | |
| 1,1-Dichloroethane | 1.648916 | 3.538465 | 10.822 | 3.205983E-02 | | | 30 | |
| 1,1-Dichloroethylene | 1.385023 | 4.865253 | 8.8476 | 6.763089E-02 | | | 30 | |
| 1,2,4-Trichlorobenzene | 0.5148378 | 41.12453 | 27.203 | 1.480613E-02 | | | | |
| 1,2,4-Trimethylbenzene | 1.051784 | 53.11966 | 22.78 | 1.048918E-02 | | | | |
| 1,2-Dibromoethane | 0.5179859 | 20.07364 | 18.371 | 0.01893 | | | 30 | |
| 1,2-Dichlorobenzene | 0.7131371 | 44.63336 | 24.4705 | 9.026326E-03 | | | | |
| 1,2-Dichloroethane | 1.069449 | 5.829632 | 13.33775 | 3.130091E-02 | | | 30 | |
| 1,2-Dichloropropane | 0.3360172 | 7.976158 | 14.7165 | 2.215717E-02 | | | 30 | |
| 1,2-Dichlorotetrafluoroethane | 2.360687 | 11.30095 | 5.649875 | 5.655805E-02 | | | 30 | |
| 1,3,5-Trimethylbenzene | 1.062278 | 43.9929 | 21.93375 | 2.119118E-02 | | | | |
| 1,3-Butadiene | 0.5735416 | 20.11683 | 6.228625 | 0.1091135 | | | 30 | |
| 1,3-Dichlorobenzene | 0.6696077 | 49.76849 | 23.59475 | 1.457982E-02 | | | | |
| 1,3-Dichloropropane | 0.4654973 | 12.69581 | 17.48538 | 2.969265E-02 | | | 30 | |
| 1,4-Dichlorobenzene | 0.6236427 | 54.56843 | 23.7785 | 0.0144024 | | | | |
| 1,4-Dioxane | 0.1846497 | 19.69241 | 15.04 | 6.478083E-02 | | | 30 | |
| 2-Butanone | 1.733964 | 44.92374 | 11.428 | 0.1635914 | | | | |
| 2-Hexanone | 0.7546478 | 6.562211 | 17.0626 | 1.673479E-02 | | | | |
| 3-Chloropropene | 1.278573 | 4.346528 | 9.42675 | 2.769754E-02 | | | 30 | |
| 4-Methyl-2-pentanone | 0.8423353 | 7.814428 | 15.6044 | 2.275779E-02 | | | | |
| Acetone | 1.47382 | 6.419353 | 8.479625 | 0.2190116 | | | 30 | |
| Acrolein | 0.4043315 | 22.66784 | 8.366 | 0.1460145 | | | 30 | |
| Acrylonitrile | 0.8090825 | 9.834437 | 9.66775 | 0.1280212 | | | 30 | |
| Benzene | 2.912755 | 10.13454 | 13.46875 | 2.178408E-02 | | | 30 | |
| Benzyl chloride | 0.2773211 | 13.16528 | 23.9242 | 9.274151E-03 | | | | |
| Bromodichloromethane | 0.5758481 | 6.5773 | 15.07712 | 6.614293E-03 | | | 30 | |
| Bromoform | 0.6537568 | 22.21779 | 20.77863 | 2.105191E-02 | | | 30 | |
| Bromomethane | 0.9395042 | 8.551225 | 7.02475 | 8.462726E-02 | | | 30 | |

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YB10001Instrument: 5975CCalibration Date: 01/28/21 00:14

| Compound | Mean RF | RF RSD | Mean RT | RT RSD | Linear r | Quad COD | LIMIT | Q |
|--------------------------------|-----------|----------|----------|--------------|----------|-----------|-------|---|
| Carbon disulfide | 2.728723 | 5.37879 | 9.72275 | 4.687565E-02 | | | 30 | |
| Carbon tetrachloride | 1.694742 | 14.68191 | 13.2706 | 4.080666E-02 | | | 30 | |
| Chlorobenzene | 0.9575035 | 7.62703 | 19.1665 | 2.056071E-02 | | | 30 | |
| Chloroethane | 0.4800299 | 7.407192 | 7.19125 | 7.785637E-02 | | | 30 | |
| Chloroform | 1.845091 | 7.059935 | 12.078 | 2.021945E-02 | | | 30 | |
| Chloromethane | 0.6590617 | 16.0044 | 5.8795 | 0.1033726 | | | 30 | |
| cis-1,2-Dichloroethylene | 1.253033 | 6.392572 | 11.8076 | 4.132915E-02 | | | 30 | |
| cis-1,3-Dichloropropylene | 0.4408728 | 11.38988 | 15.958 | 2.476346E-02 | | | 30 | |
| Cyclohexane | 1.649441 | 4.298776 | 12.97475 | 0.0192325 | | | 30 | |
| Dibromochloromethane | 0.5839562 | 15.28362 | 17.9995 | 1.987087E-02 | | | 30 | |
| Dichlorodifluoromethane | 2.322625 | 13.24178 | 5.39075 | 4.249008E-02 | | | 30 | |
| Ethanol | 0.4594539 | 12.97662 | 7.3305 | 0.2174029 | | | 30 | |
| Ethyl acetate | 1.937022 | 44.83897 | 11.72213 | 0.1082793 | | | | |
| Ethyl Benzene | 1.410567 | 10.90795 | 19.266 | 2.250315E-03 | | | 30 | |
| Hexachlorobutadiene | 0.5938052 | 47.83308 | 27.47587 | 2.302084E-02 | | | | |
| Isopropanol | 1.700061 | 30.15725 | 8.253875 | 0.2477185 | | | | |
| Isopropylbenzene | 1.382457 | 29.82945 | 20.8465 | 1.940285E-02 | | | 30 | |
| Methyl Methacrylate | 0.2362587 | 54.33968 | 14.72625 | 4.321529E-02 | | | | |
| Methyl tert-butyl ether (MTBE) | 1.975569 | 17.38184 | 9.98025 | 0.1302265 | | | 30 | |
| Methylene chloride | 2.108219 | 71.28229 | 9.580429 | 2.787056E-02 | | 0.9999807 | 0.99 | |
| Naphthalene | 1.605713 | 28.81616 | 27.6004 | 9.631682E-03 | | | | |
| n-Butylbenzene | 1.745531 | 5.239689 | 24.2052 | 1.039977E-02 | | | | |
| n-Heptane | 1.979354 | 3.533298 | 13.47325 | 1.698537E-02 | | | 30 | |
| n-Hexane | 1.509872 | 4.184026 | 10.45075 | 2.661884E-02 | | | 30 | |
| n-Propylbenzene | 1.619682 | 42.32279 | 21.62125 | 1.170543E-02 | | | | |
| o-Xylene | 1.045352 | 18.98381 | 20.19662 | 4.30727E-03 | | | 30 | |
| p- & m- Xylenes | 1.044721 | 14.21382 | 19.41275 | 1.990121E-02 | | | 30 | |
| p-Ethyltoluene | 1.257361 | 48.83878 | 21.84562 | 1.831582E-02 | | | | |
| p-Isopropyltoluene | 1.327251 | 60.94287 | 23.42 | 2.321695E-02 | | | | |
| Propylene | 0.6547622 | 10.1171 | 5.31925 | 0.084912 | | | 30 | |
| sec-Butylbenzene | 1.63111 | 52.59124 | 23.161 | 0.0131555 | | | | |
| Styrene | 0.7443945 | 35.85395 | 20.21475 | 0.0108606 | | | | |

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YB10001Instrument: 5975CCalibration Date: 01/28/21 00:14

| Compound | Mean RF | RF RSD | Mean RT | RT RSD | Linear r | Quad COD | LIMIT | Q |
|-----------------------------------|-----------|----------|----------|--------------|----------|----------|-------|---|
| SURR: p-Bromofluorobenzene | 0.5337658 | 13.1414 | 21.26575 | 6.316804E-03 | | | 30 | |
| tert-Butylbenzene | 1.19972 | 51.03046 | 22.70475 | 1.377968E-02 | | | | |
| Tetrachloroethylene | 0.4338053 | 8.675105 | 17.80325 | 1.941832E-02 | | | 30 | |
| Tetrahydrofuran | 1.034869 | 37.52092 | 12.41175 | 0.1445892 | | | | |
| Toluene | 0.9712132 | 5.892857 | 16.61325 | 1.792639E-02 | | | 30 | |
| trans-1,2-Dichloroethylene | 1.319287 | 4.934335 | 10.1745 | 4.370982E-02 | | | 30 | |
| trans-1,3-Dichloropropylene | 0.3622548 | 21.25642 | 16.7675 | 2.834501E-02 | | | 30 | |
| Trichloroethylene | 0.3906294 | 13.17685 | 14.4926 | 6.252852E-03 | | | 30 | |
| Trichlorofluoromethane (Freon 11) | 1.968664 | 10.24719 | 7.76275 | 2.933307E-02 | | | 30 | |
| Vinyl acetate | 2.038641 | 57.8008 | 10.76787 | 8.271396E-02 | | | | |
| Vinyl bromide | 0.8424028 | 10.29368 | 7.6475 | 5.681086E-02 | | | 30 | |
| Vinyl Chloride | 0.8529717 | 32.51032 | 6.1503 | 0.1429984 | | | | |

FORM VII

CONTINUING CALIBRATION CHECK

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 21A0950Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYInstrument ID: TO15_AIR2Calibration: YA10030Lab File ID: TQ213076.DCalibration Date: 01/26/21 07:48Sequence: Y1B0112Injection Date: 01/30/21Lab Sample ID: Y1B0112-CCV1Injection Time: 01:43

| COMPOUND | TYPE | CONC. (ppbv) | | RESPONSE FACTOR | | | % DIFF / DRIFT | |
|--------------------------------|------|--------------|-------|-----------------|--------------|---------|----------------|-----------|
| | | STD | CCV | ICAL | CCV | MIN (#) | CCV | LIMIT (#) |
| Bromoform | A | 10.0 | 12.0 | 0.6920348 | 0.8279185 | | 19.6 | 30 |
| Bromomethane | A | 10.0 | 9.42 | 1.185133 | 1.116616 | | -5.8 | 30 |
| Carbon disulfide | A | 10.0 | 10.9 | 3.328068 | 3.617659 | | 8.7 | 30 |
| Carbon tetrachloride | A | 10.0 | 12.1 | 3.480266 | 4.071391 | | 17.0 | 30 |
| Chlorobenzene | A | 10.0 | 10.2 | 1.052411 | 1.075185 | | 2.2 | 30 |
| Chloroethane | A | 10.0 | 9.28 | 0.4987259 | 0.4629157 | | -7.2 | 30 |
| Chloroform | A | 10.0 | 11.0 | 3.027174 | 3.340547 | | 10.4 | 30 |
| Chloromethane | A | 10.0 | 10.9 | 0.519793 | 0.5671055 | | 9.1 | 30 |
| cis-1,2-Dichloroethylene | A | 10.0 | 10.4 | 1.593994 | 1.654451 | | 3.8 | 30 |
| cis-1,3-Dichloropropylene | A | 10.0 | 10.7 | 0.4529623 | 0.4840344 | | 6.9 | 30 |
| Cyclohexane | A | 10.0 | 10.4 | 1.4958 | 1.563022 | | 4.5 | 30 |
| Dibromochloromethane | A | 10.0 | 11.4 | 0.696375 | 0.7968079 | | 14.4 | 30 |
| Dichlorodifluoromethane | A | 10.0 | 3.15 | 3.593469 | 1.133352 | | -68.5 | 30 * |
| Ethyl acetate | A | 10.0 | 10.4 | 1.934383 | 2.014633 | | 4.1 | 30 |
| Ethyl Benzene | A | 10.0 | 10.6 | 1.578363 | 1.676126 | | 6.2 | 30 |
| Hexachlorobutadiene | A | 10.0 | 12.1 | 0.7759784 | 0.9359681 | | 20.6 | 30 |
| Isopropanol | Q | 10.0 | 10.0 | 1.736529 | 1.737327 | | 0.0 | 30 |
| Methyl Methacrylate | A | 10.0 | 10.4 | 0.2618377 | 0.272681 | | 4.1 | 30 |
| Methyl tert-butyl ether (MTBE) | A | 10.0 | 10.8 | 3.649256 | 3.950155 | | 8.2 | 30 |
| Methylene chloride | A | 10.0 | 9.57 | 1.086933 | 1.040257 | | -4.3 | 30 |
| n-Heptane | A | 10.0 | 10.6 | 1.409481 | 1.497145 | | 6.2 | 30 |
| n-Hexane | A | 10.0 | 9.07 | 1.640204 | 1.487098 | | -9.3 | 30 |
| o-Xylene | A | 10.0 | 10.8 | 1.234236 | 1.330793 | | 7.8 | 30 |
| p- & m- Xylenes | A | 20.0 | 21.8 | 1.182349 | 1.289487 | | 9.1 | 30 |
| p-Ethyltoluene | A | 10.0 | 12.0 | 1.716013 | 2.053448 | | 19.7 | 30 |
| Propylene | A | 10.0 | 0.370 | 0.3825546 | 1.423775E-02 | | -96.3 | 30 * |
| Styrene | A | 10.0 | 11.6 | 0.9084209 | 1.051449 | | 15.7 | 30 |
| Tetrachloroethylene | A | 10.0 | 11.0 | 0.607179 | 0.6669628 | | 9.8 | 30 |
| Tetrahydrofuran | A | 10.0 | 10.1 | 0.9039142 | 0.9126371 | | 1.0 | 30 |

ATTACHMENT D

ANNOTATED SUMMARY FORMS

Laboratory: York Analytical Laboratories, Inc. SDG: 21A0950
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 21A0950-01 File ID: TO287826.D
 Sampled: 01/22/21 14:30 Prepared: 01/28/21 09:00 Analyzed: 01/29/21 11:48
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BA11299 Sequence: Y1A2944 Calibration: YB10001 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|----|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 23.1 | 16 | UJ |
| 71-55-6 | 1,1,1-Trichloroethane | 23.1 | 13 | UJ |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 23.1 | 16 | UJ |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 23.1 | 18 | UJ |
| 79-00-5 | 1,1,2-Trichloroethane | 23.1 | 13 | UJ |
| 75-34-3 | 1,1-Dichloroethane | 23.1 | 9.3 | UJ |
| 75-35-4 | 1,1-Dichloroethylene | 23.1 | 670 | J |
| 120-82-1 | 1,2,4-Trichlorobenzene | 23.1 | 341 | R |
| 95-63-6 | 1,2,4-Trimethylbenzene | 23.1 | 11 | UJ |
| 106-93-4 | 1,2-Dibromoethane | 23.1 | 18 | UJ |
| 95-50-1 | 1,2-Dichlorobenzene | 23.1 | 417 | UJ |
| 107-06-2 | 1,2-Dichloroethane | 23.1 | 9.3 | UJ |
| 78-87-5 | 1,2-Dichloropropane | 23.1 | 11 | UJ |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 23.1 | 16 | UJ |
| 108-67-8 | 1,3,5-Trimethylbenzene | 23.1 | 341 | UJ |
| 106-99-0 | 1,3-Butadiene | 23.1 | 15 | UJ |
| 541-73-1 | 1,3-Dichlorobenzene | 23.1 | 417 | UJ |
| 142-28-9 | 1,3-Dichloropropane | 23.1 | 11 | UJ |
| 106-46-7 | 1,4-Dichlorobenzene | 23.1 | 417 | UJ |
| 123-91-1 | 1,4-Dioxane | 23.1 | 17 | UJ |
| 78-93-3 | 2-Butanone | 23.1 | 204 | UJ |
| 591-78-6 | 2-Hexanone | 23.1 | 19 | UJ |
| 107-05-1 | 3-Chloropropene | 23.1 | 36 | UJ |
| 108-10-1 | 4-Methyl-2-pentanone | 23.1 | 9.5 | UJ |
| 67-64-1 | Acetone | 23.1 | 45 | J |
| 107-13-1 | Acrylonitrile | 23.1 | 5.0 | UJ |
| 71-43-2 | Benzene | 23.1 | 26 | J |
| 100-44-7 | Benzyl chloride | 23.1 | 12 | UJ |
| 75-27-4 | Bromodichloromethane | 23.1 | 15 | UJ |
| 75-25-2 | Bromoform | 23.1 | 24 | UJ |
| 74-83-9 | Bromomethane | 23.1 | 9.0 | UJ |
| 75-15-0 | Carbon disulfide | 23.1 | 13 | J |
| 56-23-5 | Carbon tetrachloride | 23.1 | 3.6 | UJ |
| 108-90-7 | Chlorobenzene | 23.1 | 11 | UJ |
| 75-00-3 | Chloroethane | 23.1 | 70 | J |
| 67-66-3 | Chloroform | 23.1 | 11 | J |
| 74-87-3 | Chloromethane | 23.1 | 4.8 | UJ |
| 10061-01-5 | cis-1,3-Dichloropropylene | 23.1 | 10 | UJ |
| 110-82-7 | Cyclohexane | 23.1 | 82 | UJ |
| 124-48-1 | Dibromochloromethane | 23.1 | 20 | UJ |

Laboratory: York Analytical Laboratories, Inc. SDG: 21A0950
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 21A0950-01 File ID: TO287826.D
 Sampled: 01/22/21 14:30 Prepared: 01/28/21 09:00 Analyzed: 01/29/21 11:48
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BA11299 Sequence: Y1A2944 Calibration: YB10001 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 75-71-8 | Dichlorodifluoromethane | 23.1 | 11 | UJ |
| 141-78-6 | Ethyl acetate | 23.1 | 250 | UJ |
| 100-41-4 | Ethyl Benzene | 23.1 | 10 | UJ |
| 87-68-3 | Hexachlorobutadiene | 23.1 | 25 | UJ |
| 67-63-0 | Isopropanol | 23.1 | 254 | UJ |
| 80-62-6 | Methyl Methacrylate | 23.1 | 240 | UJ |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 23.1 | 8.3 | UJ |
| 75-09-2 | Methylene chloride | 23.1 | 241 | UJ |
| 142-82-5 | n-Heptane | 23.1 | 87 | J |
| 110-54-3 | n-Hexane | 23.1 | 54 | J |
| 95-47-6 | o-Xylene | 23.1 | 10 | UJ |
| 179601-23-1 | p- & m- Xylenes | 23.1 | 20 | UJ |
| 622-96-8 | p-Ethyltoluene | 23.1 | 341 | UJ |
| 115-07-1 | Propylene | 23.1 | 4.0 | UJ |
| 100-42-5 | Styrene | 23.1 | 295 | UJ |
| 109-99-9 | Tetrahydrofuran | 23.1 | 204 | UJ |
| 108-88-3 | Toluene | 23.1 | 21 | J |
| 156-60-5 | trans-1,2-Dichloroethylene | 23.1 | 820 | J |
| 10061-02-6 | trans-1,3-Dichloropropylene | 23.1 | 10 | UJ |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 23.1 | 13 | UJ |
| 108-05-4 | Vinyl acetate | 23.1 | 244 | UJ |
| 593-60-2 | Vinyl bromide | 23.1 | 10 | UJ |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 343584 | 12.31 | 340269 | 12.304 | |
| ISTD: 1,4-Difluorobenzene | 1108308 | 13.858 | 1154282 | 13.852 | |
| ISTD: d5-Chlorobenzene | 826775 | 19.101 | 987294 | 19.095 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. SDG: 21A0950
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 21A0950-01RE1 File ID: TO287839.D
 Sampled: 01/22/21 14:30 Prepared: 01/29/21 18:00 Analyzed: 01/30/21 02:25
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BA11424 Sequence: Y1B0113 Calibration: YB10001 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|----------|----------------------------|----------|----------------------------|---|
| 127-18-4 | Tetrachloroethylene | 247 | 73000 | J |
| 79-01-6 | Trichloroethylene | 247 | 28000 | J |
| 75-01-4 | Vinyl Chloride | 247 | 7900 | J |
| 460-00-4 | SURR: p-Bromofluorobenzene | 247 | 8.1 | J |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 343582 | 12.304 | 354135 | 12.304 | |
| ISTD: 1,4-Difluorobenzene | 1046080 | 13.852 | 1201521 | 13.852 | |
| ISTD: d5-Chlorobenzene | 758481 | 19.095 | 1029256 | 19.095 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. SDG: 21A0950
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 21A0950-01RE2 File ID: TO287840.D
 Sampled: 01/22/21 14:30 Prepared: 01/29/21 18:00 Analyzed: 01/30/21 03:12
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BA11424 Sequence: Y1B0113 Calibration: YB10001 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|----------|----------------------------|----------|----------------------------|---|
| 156-59-2 | cis-1,2-Dichloroethylene | 662 | 79000 | J |
| 460-00-4 | SURR: p-Bromofluorobenzene | 662 | 8.1 | J |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|--------|--------|----------|--------|---|
| Bromochloromethane | 323860 | 12.304 | 354135 | 12.304 | |
| ISTD: 1,4-Difluorobenzene | 985953 | 13.852 | 1201521 | 13.852 | |
| ISTD: d5-Chlorobenzene | 686290 | 19.101 | 1029256 | 19.095 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. SDG: 21A0950
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 21A0950-02 File ID: TO287841.D
 Sampled: 01/22/21 14:30 Prepared: 01/29/21 18:00 Analyzed: 01/30/21 04:04
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BA11424 Sequence: Y1B0113 Calibration: YB10001 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|----|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.69 | 1.2 | UJ |
| 71-55-6 | 1,1,1-Trichloroethane | 1.69 | 0.92 | UJ |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.69 | 1.2 | UJ |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 1.69 | 1.3 | UJ |
| 79-00-5 | 1,1,2-Trichloroethane | 1.69 | 0.92 | UJ |
| 75-34-3 | 1,1-Dichloroethane | 1.69 | 0.68 | UJ |
| 75-35-4 | 1,1-Dichloroethylene | 1.69 | 0.17 | UJ |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.69 | 1.3 | R |
| 95-63-6 | 1,2,4-Trimethylbenzene | 1.69 | 1.4 | J |
| 106-93-4 | 1,2-Dibromoethane | 1.69 | 1.3 | UJ |
| 95-50-1 | 1,2-Dichlorobenzene | 1.69 | 30.5 | UJ |
| 107-06-2 | 1,2-Dichloroethane | 1.69 | 0.68 | UJ |
| 78-87-5 | 1,2-Dichloropropane | 1.69 | 0.78 | UJ |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 1.69 | 1.2 | UJ |
| 108-67-8 | 1,3,5-Trimethylbenzene | 1.69 | 24.9 | UJ |
| 106-99-0 | 1,3-Butadiene | 1.69 | 1.1 | UJ |
| 541-73-1 | 1,3-Dichlorobenzene | 1.69 | 30.5 | UJ |
| 142-28-9 | 1,3-Dichloropropane | 1.69 | 0.78 | UJ |
| 106-46-7 | 1,4-Dichlorobenzene | 1.69 | 30.5 | UJ |
| 123-91-1 | 1,4-Dioxane | 1.69 | 1.2 | UJ |
| 78-93-3 | 2-Butanone | 1.69 | 16 | J |
| 591-78-6 | 2-Hexanone | 1.69 | 3.2 | J |
| 107-05-1 | 3-Chloropropene | 1.69 | 2.6 | UJ |
| 108-10-1 | 4-Methyl-2-pentanone | 1.69 | 1.7 | J |
| 107-13-1 | Acrylonitrile | 1.69 | 0.37 | UJ |
| 71-43-2 | Benzene | 1.69 | 2.7 | J |
| 100-44-7 | Benzyl chloride | 1.69 | 0.87 | UJ |
| 75-27-4 | Bromodichloromethane | 1.69 | 1.1 | UJ |
| 75-25-2 | Bromoform | 1.69 | 1.7 | UJ |
| 74-83-9 | Bromomethane | 1.69 | 0.66 | UJ |
| 75-15-0 | Carbon disulfide | 1.69 | 8.3 | J |
| 56-23-5 | Carbon tetrachloride | 1.69 | 0.32 | J |
| 108-90-7 | Chlorobenzene | 1.69 | 0.78 | UJ |
| 75-00-3 | Chloroethane | 1.69 | 0.45 | UJ |
| 67-66-3 | Chloroform | 1.69 | 2.4 | J |
| 74-87-3 | Chloromethane | 1.69 | 1.4 | J |
| 156-59-2 | cis-1,2-Dichloroethylene | 1.69 | 65 | J |
| 10061-01-5 | cis-1,3-Dichloropropylene | 1.69 | 0.77 | UJ |
| 110-82-7 | Cyclohexane | 1.69 | 1.2 | J |
| 124-48-1 | Dibromochloromethane | 1.69 | 1.4 | UJ |

Laboratory: York Analytical Laboratories, Inc. SDG: 21A0950
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 21A0950-02 File ID: TO287841.D
 Sampled: 01/22/21 14:30 Prepared: 01/29/21 18:00 Analyzed: 01/30/21 04:04
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BA11424 Sequence: Y1B0113 Calibration: YB10001 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 75-71-8 | Dichlorodifluoromethane | 1.69 | 2.0 | J |
| 141-78-6 | Ethyl acetate | 1.69 | 18.3 | UJ |
| 100-41-4 | Ethyl Benzene | 1.69 | 1.3 | J |
| 87-68-3 | Hexachlorobutadiene | 1.69 | 1.8 | UJ |
| 67-63-0 | Isopropanol | 1.69 | 1.7 | J |
| 80-62-6 | Methyl Methacrylate | 1.69 | 20.8 | UJ |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 1.69 | 0.97 | J |
| 75-09-2 | Methylene chloride | 1.69 | 5.7 | J |
| 142-82-5 | n-Heptane | 1.69 | 61 | J |
| 110-54-3 | n-Hexane | 1.69 | 21 | J |
| 95-47-6 | o-Xylene | 1.69 | 1.6 | J |
| 179601-23-1 | p- & m- Xylenes | 1.69 | 4.3 | J |
| 622-96-8 | p-Ethyltoluene | 1.69 | 1.4 | J |
| 115-07-1 | Propylene | 1.69 | 4.4 | J |
| 100-42-5 | Styrene | 1.69 | 21.6 | UJ |
| 127-18-4 | Tetrachloroethylene | 1.69 | 400 | UJ |
| 109-99-9 | Tetrahydrofuran | 1.69 | 2.8 | J |
| 108-88-3 | Toluene | 1.69 | 7.8 | J |
| 156-60-5 | trans-1,2-Dichloroethylene | 1.69 | 0.67 | UJ |
| 10061-02-6 | trans-1,3-Dichloropropylene | 1.69 | 0.77 | UJ |
| 79-01-6 | Trichloroethylene | 1.69 | 26 | J |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 1.69 | 1.0 | J |
| 108-05-4 | Vinyl acetate | 1.69 | 17.8 | UJ |
| 593-60-2 | Vinyl bromide | 1.69 | 0.74 | UJ |
| 75-01-4 | Vinyl Chloride | 1.69 | 1.9 | J |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 326120 | 12.304 | 354135 | 12.304 | |
| ISTD: 1,4-Difluorobenzene | 1074897 | 13.852 | 1201521 | 13.852 | |
| ISTD: d5-Chlorobenzene | 963577 | 19.095 | 1029256 | 19.095 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. SDG: 21A0950
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 21A0950-02RE1 File ID: TO287829.D
 Sampled: 01/22/21 14:30 Prepared: 01/28/21 09:00 Analyzed: 01/29/21 14:21
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BA11299 Sequence: Y1A2944 Calibration: YB10001 Instrument: 5975C

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|----------|----------------------------|----------|----------------------------|---|
| 67-64-1 | Acetone | 16.9 | 190 | J |
| 460-00-4 | SURR: p-Bromofluorobenzene | 16.9 | 9.7 | J |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 406177 | 12.304 | 340269 | 12.304 | |
| ISTD: 1,4-Difluorobenzene | 1333166 | 13.852 | 1154282 | 13.852 | |
| ISTD: d5-Chlorobenzene | 1017478 | 19.095 | 987294 | 19.095 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. SDG: 21A0950
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Indoor Ambient Air Laboratory ID: 21A0950-03 File ID: TQ213084.D
 Sampled: 01/22/21 16:00 Prepared: 01/29/21 18:00 Analyzed: 01/30/21 09:18
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BA11425 Sequence: Y1B0112 Calibration: YA10030 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|----|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.04 | 0.71 | UJ |
| 71-55-6 | 1,1,1-Trichloroethane | 1.04 | 0.57 | UJ |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.04 | 0.71 | UJ |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 1.04 | 0.79 | UJ |
| 79-00-5 | 1,1,2-Trichloroethane | 1.04 | 0.57 | UJ |
| 75-34-3 | 1,1-Dichloroethane | 1.04 | 0.42 | UJ |
| 75-35-4 | 1,1-Dichloroethylene | 1.04 | 0.10 | UJ |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.04 | 0.77 | R |
| 95-63-6 | 1,2,4-Trimethylbenzene | 1.04 | 1.0 | J |
| 106-93-4 | 1,2-Dibromoethane | 1.04 | 0.80 | UJ |
| 95-50-1 | 1,2-Dichlorobenzene | 1.04 | 0.62 | UJ |
| 107-06-2 | 1,2-Dichloroethane | 1.04 | 0.42 | UJ |
| 78-87-5 | 1,2-Dichloropropane | 1.04 | 0.48 | UJ |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 1.04 | 0.72 | UJ |
| 108-67-8 | 1,3,5-Trimethylbenzene | 1.04 | 0.51 | UJ |
| 106-99-0 | 1,3-Butadiene | 1.04 | 0.69 | UJ |
| 541-73-1 | 1,3-Dichlorobenzene | 1.04 | 17.3 | UJ |
| 142-28-9 | 1,3-Dichloropropane | 1.04 | 0.48 | UJ |
| 106-46-7 | 1,4-Dichlorobenzene | 1.04 | 17.3 | UJ |
| 123-91-1 | 1,4-Dioxane | 1.04 | 0.75 | UJ |
| 78-93-3 | 2-Butanone | 1.04 | 2.9 | J |
| 591-78-6 | 2-Hexanone | 1.04 | 0.85 | UJ |
| 107-05-1 | 3-Chloropropene | 1.04 | 1.6 | UJ |
| 108-10-1 | 4-Methyl-2-pentanone | 1.04 | 0.42 | UJ |
| 67-64-1 | Acetone | 1.04 | 12 | J |
| 107-13-1 | Acrylonitrile | 1.04 | 0.23 | UJ |
| 71-43-2 | Benzene | 1.04 | 1.1 | J |
| 100-44-7 | Benzyl chloride | 1.04 | 15.5 | UJ |
| 75-27-4 | Bromodichloromethane | 1.04 | 0.69 | UJ |
| 75-25-2 | Bromoform | 1.04 | 1.1 | UJ |
| 74-83-9 | Bromomethane | 1.04 | 0.40 | UJ |
| 75-15-0 | Carbon disulfide | 1.04 | 0.32 | UJ |
| 56-23-5 | Carbon tetrachloride | 1.04 | 0.52 | J |
| 108-90-7 | Chlorobenzene | 1.04 | 0.48 | UJ |
| 75-00-3 | Chloroethane | 1.04 | 0.27 | UJ |
| 67-66-3 | Chloroform | 1.04 | 0.56 | J |
| 74-87-3 | Chloromethane | 1.04 | 1.2 | J |
| 156-59-2 | cis-1,2-Dichloroethylene | 1.04 | 35 | J |
| 10061-01-5 | cis-1,3-Dichloropropylene | 1.04 | 0.47 | UJ |
| 110-82-7 | Cyclohexane | 1.04 | 0.57 | J |

Laboratory: York Analytical Laboratories, Inc. SDG: 21A0950
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Indoor Ambient Air Laboratory ID: 21A0950-03 File ID: TQ213084.D
 Sampled: 01/22/21 16:00 Prepared: 01/29/21 18:00 Analyzed: 01/30/21 09:18
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BA11425 Sequence: Y1B0112 Calibration: YA10030 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 124-48-1 | Dibromochloromethane | 1.04 | 0.88 | UJ |
| 75-71-8 | Dichlorodifluoromethane | 1.04 | 1.6 | J |
| 141-78-6 | Ethyl acetate | 1.04 | 0.75 | UJ |
| 100-41-4 | Ethyl Benzene | 1.04 | 0.63 | J |
| 87-68-3 | Hexachlorobutadiene | 1.04 | 1.1 | UJ |
| 67-63-0 | Isopropanol | 1.04 | 17 | J |
| 80-62-6 | Methyl Methacrylate | 1.04 | 1.4 | J |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 1.04 | 0.37 | UJ |
| 75-09-2 | Methylene chloride | 1.04 | 2.6 | J |
| 142-82-5 | n-Heptane | 1.04 | 2.8 | J |
| 110-54-3 | n-Hexane | 1.04 | 1.8 | J |
| 95-47-6 | o-Xylene | 1.04 | 0.86 | J |
| 179601-23-1 | p- & m- Xylenes | 1.04 | 2.2 | J |
| 622-96-8 | p-Ethyltoluene | 1.04 | 0.82 | J |
| 115-07-1 | Propylene | 1.04 | 0.18 | R |
| 100-42-5 | Styrene | 1.04 | 0.44 | UJ |
| 127-18-4 | Tetrachloroethylene | 1.04 | 19 | J |
| 109-99-9 | Tetrahydrofuran | 1.04 | 3.4 | J |
| 108-88-3 | Toluene | 1.04 | 3.8 | J |
| 156-60-5 | trans-1,2-Dichloroethylene | 1.04 | 0.41 | UJ |
| 10061-02-6 | trans-1,3-Dichloropropylene | 1.04 | 0.47 | UJ |
| 79-01-6 | Trichloroethylene | 1.04 | 7.2 | J |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 1.04 | 2.1 | J |
| 108-05-4 | Vinyl acetate | 1.04 | 0.37 | UJ |
| 593-60-2 | Vinyl bromide | 1.04 | 0.45 | UJ |
| 75-01-4 | Vinyl Chloride | 1.04 | 1.0 | J |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 288825 | 12.036 | 232902 | 12.039 | |
| ISTD: 1,4-Difluorobenzene | 1494646 | 13.618 | 1292063 | 13.621 | |
| ISTD: d5-Chlorobenzene | 1271520 | 18.884 | 1203604 | 18.884 | |

* Values outside of QC limits

JR 12/8/2025

December 8, 2025

Mr. Jason Stewart
Advanced Cleanup Technologies
228 Park Ave S PMB 34864
New York, New York 10003

Re: Data Usability Summary Report – York Analytical Laboratories, Inc. - 21D0660

Dear Mr. Stewart:

The evaluation of analytical data by York Analytical Laboratories for project 9268-BXNY, which were reported in a single data package under Job No. 21D0660 has been completed. The following samples were reported.

SV-1 SV-2 SV-3

Analysis was performed in accordance with EPA Method TO-15 (volatile organics). The review was performed to the extent possible, in accordance with the analytical method, and “DER-10/ Technical Guidance for Site Investigation and Remediation”. Professional judgment is applied as necessary and appropriate. National Functional Guidelines for Organic Data Review was consulted as needed. Qualifiers consistent with those defined by EPA Region 2 are applied as necessary and appropriate.

Below is the Data Usability Summary Report (DUSR) associated with these samples.

| Data Usability Summary Report | |
|---|----------------------------|
| 1. Is the data package complete as defined under the requirements for the most current DEC ASP Category B or USEPA CLP data deliverables? | Yes |
| 2. Have all holding times been met? | Yes |
| 3. Do all the QC data; blanks, instrument tunings, calibration standards, calibration verifications, surrogate recoveries, spike recoveries, replicate analyses, laboratory controls and sample data fall within the protocol required limits and specifications? | No -see following sections |
| 4. Have all of the data been generated using established and agreed upon analytical protocols? | Yes |
| 5. Does an evaluation of the raw data confirm the results provided in the data summary sheet and the quality control verification forms? | Yes |
| 6. Have the correct data qualifiers been used and are they consistent with the most current DEC ASP? | Yes |
| 7. Have any quality control (QC) exceedances been specifically noted in the DUSR and have the corresponding QC summary sheet from the data package been attached to the DUSR? | Yes |

Overall Evaluation

Based on the data review effort, results are usable, with the following qualifications. For samples that are qualified as estimated (J-, UJ), detected results may be biased low. False negatives may exist in non-detect results. Sample results that are qualified as estimated (J+) may be biased high. For samples that are qualified as estimated with any combination of (J), (J-) and/or (J+), the (J) qualifier takes precedence and is applied to the sample result. It is not possible to determine the direction of the bias and the overall effect on the result.

- The result for methylene chloride in SV-1 and SV-2 is qualified as estimated (J+) due to method blank contamination.
- The result for methylene chloride in SV-1 and SV-2 is qualified as estimated (J+) due to detection in the canister used for clean canister certification.

Qualifier definitions are provided in Attachment A. A copy of the chain of custody record is provided in Attachment B. Pages from the data package illustrating the exceedances and issues described in this validation report are included in Attachment C. Annotated Summary Forms are included in Attachment D detailing qualifications resulting from the data review effort.

The following components were reviewed, where applicable:

- Chain of Custody
- Receiving conditions
- Holding times
- Preservation
- Analyte lists
- Reporting limits
- Requested methods
- Units, and
- Sample related quality control data:
 - Method, instrument blanks
 - Clean canister certification
 - Field blanks
 - Surrogate recoveries
 - LCS recoveries
 - Internal standard area response
 - Duplicates
 - Analyte Identification
- Instrument related quality control data:
 - Instrument tunes
 - Calibration summaries

The following sections of the report detail only quality control exceedances that impacted results. Where a quality control item exceeded control limits but there is no impact to the samples results, these are not detailed in the report.

Documentation: A completeness review of the data package was performed, and the data package was determined to be a complete Category B data package. The following documentation issues were observed during the review:

- The data package indicates that the laboratory did not hold certification for ethyl acetate, p-ethyltoluene, propylene, tetrahydrofuran, 1,1,1,2-tetrachloroethane, 1,3-dichloropropane, and 2-hexanone. It is noted that NY ELAP does not offer certification for ethyl acetate, p-ethyltoluene, propylene, tetrahydrofuran and 2-hexanone, but does offer certification for 1,1,1,2-tetrachloroethane and 1,3-dichloropropane. The laboratory was contacted to request clarification. The laboratory stated *'Given the age of the project, it is difficult to ascertain the exact circumstances in which these compounds were requested. We don't typically report non-certified compounds, except at the request of the client, so it can be safely assumed that there was correspondence regarding the accreditation status of the compounds at the time the samples were taken. Just to note- while these compounds are indeed non-certified under our scope of accreditation between 2020-2022, we ran the compounds in the exact same way we ran our certified compounds – quantitated against a multiple calibration, where they were properly qualified as specified by our certifying body.'*

Holding Times, Preservation, Sample Integrity:

A copy of the applicable chain of custody (COC) record was included in the data package, documenting sample collection date of April 13, 2021. The samples were received at the laboratory on the same day on April 14, 2021. The samples were analyzed within method hold time. However, the data package does contain sample receipt information regarding condition of canisters upon receipt. The sample preparation bench sheets documents canister vacuum upon receipt, which indicates all canisters preserved (pressure inside the canister maintained within +/- 5 psi from sampling to check in the laboratory or analysis).

A. Volatile Organics

1. Blanks

The table below summarizes method blank detections that impacted sample results:

| Blank | Analyte | Conc (on-column ppbv_ | Affected Sample |
|--------------|--------------------|-----------------------|-----------------|
| BD11066-BLK1 | Methylene chloride | 0.13 | SV-2 |
| BD11103-BLK1 | | 0.13 | SV-1 |

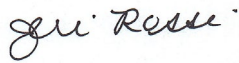
The result for methylene chloride in SV-1 and SV-2 is qualified as estimated (J+) due to method blank contamination.

2. Clean Canister Certification

Batch canister certification was performed on a canister that was not used to collect samples. Methylene chloride (0.11 ppbv on-column) is detected in the canister used for clean canister certification. The result for methylene chloride in SV-1 and SV-2 is qualified as estimated (J+) due to detection in the canister used for clean canister certification.

No other sample results are qualified. Please feel free to contact me at (908) 370-3431 or richjერიrossi513@gmail.com if you have any questions regarding this data package review report or need further information.

Sincerely,

A handwritten signature in cursive script that reads "Jeri Rossi".

Jeri L Rossi, CEAC

Environmental Consulting Chemist

ATTACHMENT A

Qualifier Definitions

EPA Qualifier Definitions

- U The analyte was analyzed for but was not detected above the level of the reported sample quantitation limit.
- J The result is an estimated quantity. The associated numerical value is the approximate concentration of the analyte in the sample.
- J+ The result is an estimated quantity, but the result may be biased high.
- J- The result is an estimated quantity, but the result may be biased low.
- NJ The analyte has been “tentatively identified” or “presumptively” as present and the associated numerical value is the estimated concentration in the sample.
- UJ The analyte was analyzed for but was not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
- R The data are unusable. The sample results are rejected due to serious deficiencies in meeting QC criteria. The analyte may or may not be present in the sample.

ATTACHMENT B

CHAIN OF CUSTODY (COC)



York Analytical Laboratories, Inc.
120 Research Drive 132-02 89th Ave Queens,
Stratford, CT 06615 NY 11418

YORK
ANALYTICAL LABORATORIES INC.

clientservices@yorklab.com
www.yorklab.com

Field Chain-of-Custody Record - AIR

YORK Project No.

2110600

NOTE: YORK's Standard Terms & Conditions are listed on the back side of this document.
This document serves as your written authorization for YORK to proceed with the analyses requested below.
signature binds you to YORK's Standard Terms & Conditions.

Your Page 1 of 1

| | | | | | | | | | |
|---|--|--------------------------------|--|---------------------------------|--|----------------------------|--|--|--|
| YOUR Information | | Report To: | | Invoice To: | | YOUR Project Number | | Turn-Around Time | |
| Company: <u>Advanced CleanTech</u> | | Company: <u>ACT</u> | | Company: <u>ACT</u> | | <u>9628-BXNY</u> | | RUSH - Next Day | |
| Address: <u>250 Brookhaven Rd Newville NY</u> | | Address: <u>same</u> | | Address: <u>same</u> | | | | RUSH - Two Day | |
| Phone.: <u>516-441-5800</u> | | Phone.: _____ | | Phone.: _____ | | YOUR Project Name | | RUSH - Three Day | |
| Contact: <u>Jason Stewart</u> | | Contact: <u>Paul Stewart</u> | | Contact: <u>Kern Friedman</u> | | | | RUSH - Four Day | |
| E-mail: <u>jasons@act-earth</u> | | E-mail: <u>pauls@act-earth</u> | | E-mail: <u>KernFO@act-earth</u> | | YOUR PO#: | | Standard (5-7 Day) <input checked="" type="checkbox"/> | |

Please print clearly and legibly. All information must be complete. Samples will not be logged in and the turn-around-time clock will not begin until any questions by YORK are resolved.

Tim Yang
Samples Collected by: (print your name above and sign below)

| | | | | | | |
|--|---------------------|--|--|--|--|---|
| Air Matrix Codes AI - Indoor Ambient Air AO - Outdoor Amb. Air AE - Vapor Extraction Well/ Process Gas/Effluent <input checked="" type="checkbox"/> AS - Soil Vapor/Sub-Slab | Samples From | | Report / EDD Type (circle selections) | | | YORK Reg. Comp. Compared to the following Regulation(s): (please fill in) |
| | New York | <input checked="" type="checkbox"/> Summary Report | CT RCP | <input checked="" type="checkbox"/> Standard Excel EDD | | |
| | New Jersey | <input type="checkbox"/> QA Report | CT RCP DQA/DUE | EQUIS (Standard) | | |
| | Connecticut | <input type="checkbox"/> NY ASP A Package | NJDEP Reduced Deliv. | NYSDEC EQUIS | | |
| | Pennsylvania | <input type="checkbox"/> NY ASP B Package | NJDKQP | NJDEP SRP HazSite | | |
| Other | Other: _____ | | | | | |

Certified Canisters: Batch ___ Individual ___

Please enter the following REQUIRED Field Data

Reporting Units: ug/m³ ___ ppbv ___ ppmv ___

| Sample Identification | Date/Time Sampled | Air Matrix | Canister Vacuum Before Sampling (in Hg) | Canister Vacuum After Sampling (in Hg) | Canister ID | Flow Cont. ID | Analysis Requested |
|-----------------------|-------------------|------------|---|--|-------------|---------------|--------------------|
| SV-1 | 4/13/21 10:3 | AS | -30 | -8 | 23992 | 6867 | TO-15 |
| SV-2 | " " | " | -30 | -8 | 468 | 7419 | " |
| SV-3 | " " | " | -30 | -8 | 37004 | 7091 | " |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |

Comments:

| | | |
|--|--|------------|
| Detection Limits Required ≤ 1 ug/m ³ ___ NYSDEC V1 Limits ___ Routine Survey ___ Other ___ | Sampling Media | |
| | 6 Liter Canister <input checked="" type="checkbox"/> | Tedlar Bag |

| | | | | | |
|-----------------------------------|-----------|-----------------------------------|----------------------|-----------------------------------|----------------------|
| Samples Relinquished by / Company | Date/Time | Samples Received by / Company | Date/Time | Samples Relinquished by / Company | Date/Time |
| <u>T-York ACT</u> | | <u>YORK</u> | <u>4/14/21 12:25</u> | <u>YORK</u> | |
| Samples Received by / Company | Date/Time | Samples Relinquished by / Company | Date/Time | Samples Received by / Company | Date/Time |
| | | | | | |
| Samples Relinquished by / Company | Date/Time | Samples Received by / Company | Date/Time | Samples Received in LAB by | Date/Time |
| | | | | <u>Stewart</u> | <u>4/14/21 14:08</u> |

Page 32 of 305

ATTACHMENT C

**SELECTED PAGES FROM DATA PACKAGE –
QC EXCEEDANCES AND VALIDATION ISSUES**

FORM I

METHOD BLANK DATA SHEET
EPA TO-15

Laboratory: York Analytical Laboratories, Inc. SDG: 21D0660
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Air Laboratory ID: BD11066-BLK1 File ID: TQ214395.D
 Prepared: 04/20/21 12:32 Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Analyzed: 04/20/21 16:34 Instrument: TO15_AIR2
 Batch: BD11066 Sequence: Y1D2135 Calibration: YC10005

| CAS NO. | COMPOUND | CONC. (ug/m ³) | Q |
|-------------|--------------------------------|----------------------------|---|
| 74-83-9 | Bromomethane | 0.39 | U |
| 75-15-0 | Carbon disulfide | 0.31 | U |
| 56-23-5 | Carbon tetrachloride | 0.16 | U |
| 108-90-7 | Chlorobenzene | 0.46 | U |
| 75-00-3 | Chloroethane | 0.26 | U |
| 67-66-3 | Chloroform | 0.49 | U |
| 74-87-3 | Chloromethane | 0.21 | U |
| 156-59-2 | cis-1,2-Dichloroethylene | 0.099 | U |
| 10061-01-5 | cis-1,3-Dichloropropylene | 0.45 | U |
| 110-82-7 | Cyclohexane | 0.34 | U |
| 124-48-1 | Dibromochloromethane | 0.85 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.49 | U |
| 141-78-6 | Ethyl acetate | 0.72 | U |
| 100-41-4 | Ethyl Benzene | 0.43 | U |
| 87-68-3 | Hexachlorobutadiene | 1.1 | U |
| 67-63-0 | Isopropanol | 0.49 | U |
| 80-62-6 | Methyl Methacrylate | 0.41 | U |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 0.36 | U |
| 75-09-2 | Methylene chloride | 0.69 | J |
| 142-82-5 | n-Heptane | 0.41 | U |
| 110-54-3 | n-Hexane | 0.35 | U |
| 95-47-6 | o-Xylene | 0.43 | U |
| 179601-23-1 | p- & m- Xylenes | 0.87 | U |
| 622-96-8 | p-Ethyltoluene | 0.49 | U |
| 115-07-1 | Propylene | 0.17 | U |
| 100-42-5 | Styrene | 0.43 | U |
| 127-18-4 | Tetrachloroethylene | 0.68 | U |
| 109-99-9 | Tetrahydrofuran | 0.59 | U |
| 108-88-3 | Toluene | 0.38 | U |
| 156-60-5 | trans-1,2-Dichloroethylene | 0.40 | U |

FORM I

METHOD BLANK DATA SHEET
EPA TO-15

Laboratory: York Analytical Laboratories, Inc. SDG: 21D0660
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Air Laboratory ID: BD11103-BLK1 File ID: TQ214421.D
 Prepared: 04/21/21 12:41 Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Analyzed: 04/21/21 19:41 Instrument: TO15_AIR2
 Batch: BD11103 Sequence: Y1D2220 Calibration: YC10005

| CAS NO. | COMPOUND | CONC. (ug/m ³) | Q |
|-------------|--------------------------------|----------------------------|---|
| 74-83-9 | Bromomethane | 0.39 | U |
| 75-15-0 | Carbon disulfide | 0.31 | U |
| 56-23-5 | Carbon tetrachloride | 0.16 | U |
| 108-90-7 | Chlorobenzene | 0.46 | U |
| 75-00-3 | Chloroethane | 0.26 | U |
| 67-66-3 | Chloroform | 0.49 | U |
| 74-87-3 | Chloromethane | 0.21 | U |
| 156-59-2 | cis-1,2-Dichloroethylene | 0.099 | U |
| 10061-01-5 | cis-1,3-Dichloropropylene | 0.45 | U |
| 110-82-7 | Cyclohexane | 0.34 | U |
| 124-48-1 | Dibromochloromethane | 0.85 | U |
| 75-71-8 | Dichlorodifluoromethane | 0.49 | U |
| 141-78-6 | Ethyl acetate | 0.72 | U |
| 100-41-4 | Ethyl Benzene | 0.43 | U |
| 87-68-3 | Hexachlorobutadiene | 1.1 | U |
| 67-63-0 | Isopropanol | 0.49 | U |
| 80-62-6 | Methyl Methacrylate | 0.41 | U |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 0.36 | U |
| 75-09-2 | Methylene chloride | 0.69 | J |
| 142-82-5 | n-Heptane | 0.41 | U |
| 110-54-3 | n-Hexane | 0.35 | J |
| 95-47-6 | o-Xylene | 0.43 | J |
| 179601-23-1 | p- & m- Xylenes | 0.87 | U |
| 622-96-8 | p-Ethyltoluene | 0.49 | U |
| 115-07-1 | Propylene | 0.17 | U |
| 100-42-5 | Styrene | 0.43 | U |
| 127-18-4 | Tetrachloroethylene | 0.68 | U |
| 109-99-9 | Tetrahydrofuran | 0.59 | U |
| 108-88-3 | Toluene | 0.38 | U |
| 156-60-5 | trans-1,2-Dichloroethylene | 0.40 | U |

Data Path : D:\040721\
 Data File : TO288688A.D
 Acq On : 7 Apr 2021 11:15 pm
 Operator : AS
 Sample : BD10393-BLK2
 Misc : QBTO1040721A Batch Cert. Can#37790 QBS040221A
 ALS Vial : 92 Sample Multiplier: 1
 InstName : 5975C

Quant Time: Apr 08 13:46:11 2021
 Quant Method : C:\msdchem\1\methods\AIR132.M
 Quant Title : TO15 VOC Analysis
 QLast Update : Mon Apr 05 17:06:27 2021
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|------------------------------|--------|----------------|----------|-------|--------|----------|
| ----- | | | | | | |
| Internal Standards | | | | | | |
| 1) Methane, bromochloro- | 12.243 | 49 | 418040 | 10.00 | ppbv | 0.00 |
| 38) 1,4-Difluorobenzene | 13.791 | 114 | 958406 | 10.00 | ppbv | 0.00 |
| 54) d5-Chlorobenzene | 19.034 | 117 | 654058 | 10.00 | ppbv | 0.00 |
| System Monitoring Compounds | | | | | | |
| 65) p-Bromofluorobenzene | 21.198 | 95 | 324506 | 7.78 | ppbv | 0.00 |
| Spiked Amount | 10.000 | Range 70 - 130 | Recovery | = | 77.80% | |
| Target Compounds | | | | | | |
| | | | | | | Qvalue |
| 2) Propylene | 5.275 | 42 | 849 | 0.04 | ppbv # | 25 |
| 4) 1,2-Dichlorotetrafluor... | 5.610 | 85 | 2491 | 0.02 | ppbv # | 87 |
| 15) Acetone | 8.487 | 43 | 1240 | 0.02 | ppbv # | 43 |
| 19) Methylene Chloride | 9.518 | 49 | 5486 | 0.11 | ppbv | 91 |
| 24) Hexane | 10.395 | 57 | 5409 | 0.08 | ppbv | 89 |
| 58) p- & m-Xylenes | 19.351 | 91 | 3875 | 0.04 | ppbv # | 69 |
| ----- | | | | | | |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

ATTACHMENT D

ANNOTATED SUMMARY FORMS

Laboratory: York Analytical Laboratories, Inc. SDG: 21D0660
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 21D0660-01 File ID: TQ214422.D
 Sampled: 04/13/21 15:00 Prepared: 04/21/21 12:41 Analyzed: 04/21/21 20:44
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BD11103 Sequence: Y1D2220 Calibration: YC10005 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.55 | 1.1 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.55 | 0.84 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.55 | 1.1 | U |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 1.55 | 1.2 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.55 | 0.84 | U |
| 75-34-3 | 1,1-Dichloroethane | 1.55 | 0.63 | U |
| 75-35-4 | 1,1-Dichloroethylene | 1.55 | 0.15 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.55 | 1.1 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 1.55 | 3.8 | |
| 106-93-4 | 1,2-Dibromoethane | 1.55 | 1.2 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.55 | 0.93 | U |
| 107-06-2 | 1,2-Dichloroethane | 1.55 | 0.63 | U |
| 78-87-5 | 1,2-Dichloropropane | 1.55 | 0.71 | U |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 1.55 | 1.1 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 1.55 | 1.3 | |
| 106-99-0 | 1,3-Butadiene | 1.55 | 1.0 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.55 | 0.93 | U |
| 142-28-9 | 1,3-Dichloropropane | 1.55 | 0.71 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.55 | 0.93 | U |
| 123-91-1 | 1,4-Dioxane | 1.55 | 1.1 | U |
| 78-93-3 | 2-Butanone | 1.55 | 1.3 | |
| 591-78-6 | 2-Hexanone | 1.55 | 1.3 | U |
| 107-05-1 | 3-Chloropropene | 1.55 | 2.4 | U |
| 108-10-1 | 4-Methyl-2-pentanone | 1.55 | 0.63 | U |
| 67-64-1 | Acetone | 1.55 | 11 | |
| 107-13-1 | Acrylonitrile | 1.55 | 0.34 | U |
| 71-43-2 | Benzene | 1.55 | 1.3 | |
| 100-44-7 | Benzyl chloride | 1.55 | 0.80 | U |
| 75-27-4 | Bromodichloromethane | 1.55 | 1.0 | U |
| 75-25-2 | Bromoform | 1.55 | 1.6 | U |
| 74-83-9 | Bromomethane | 1.55 | 0.60 | U |
| 75-15-0 | Carbon disulfide | 1.55 | 2.1 | |
| 56-23-5 | Carbon tetrachloride | 1.55 | 0.49 | |
| 108-90-7 | Chlorobenzene | 1.55 | 0.71 | U |
| 75-00-3 | Chloroethane | 1.55 | 0.41 | U |
| 67-66-3 | Chloroform | 1.55 | 8.2 | |
| 74-87-3 | Chloromethane | 1.55 | 0.32 | U |
| 156-59-2 | cis-1,2-Dichloroethylene | 1.55 | 0.15 | U |
| 10061-01-5 | cis-1,3-Dichloropropylene | 1.55 | 0.70 | U |
| 110-82-7 | Cyclohexane | 1.55 | 0.53 | U |

Laboratory: York Analytical Laboratories, Inc. SDG: 21D0660
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 21D0660-01 File ID: TQ214422.D
 Sampled: 04/13/21 15:00 Prepared: 04/21/21 12:41 Analyzed: 04/21/21 20:44
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BD11103 Sequence: Y1D2220 Calibration: YC10005 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 124-48-1 | Dibromochloromethane | 1.55 | 1.3 | U |
| 75-71-8 | Dichlorodifluoromethane | 1.55 | 2.3 | |
| 141-78-6 | Ethyl acetate | 1.55 | 1.1 | U |
| 100-41-4 | Ethyl Benzene | 1.55 | 10 | |
| 87-68-3 | Hexachlorobutadiene | 1.55 | 1.6 | U |
| 67-63-0 | Isopropanol | 1.55 | 0.76 | U |
| 80-62-6 | Methyl Methacrylate | 1.55 | 0.63 | U |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 1.55 | 0.56 | U |
| 75-09-2 | Methylene chloride | 1.55 | 1.1 | J+ |
| 142-82-5 | n-Heptane | 1.55 | 0.63 | U |
| 110-54-3 | n-Hexane | 1.55 | 0.93 | |
| 95-47-6 | o-Xylene | 1.55 | 7.0 | |
| 179601-23-1 | p- & m- Xylenes | 1.55 | 22 | |
| 622-96-8 | p-Ethyltoluene | 1.55 | 3.0 | |
| 115-07-1 | Propylene | 1.55 | 2.8 | |
| 100-42-5 | Styrene | 1.55 | 0.66 | U |
| 127-18-4 | Tetrachloroethylene | 1.55 | 95 | |
| 109-99-9 | Tetrahydrofuran | 1.55 | 1.1 | |
| 108-88-3 | Toluene | 1.55 | 17 | |
| 156-60-5 | trans-1,2-Dichloroethylene | 1.55 | 0.61 | U |
| 10061-02-6 | trans-1,3-Dichloropropylene | 1.55 | 0.70 | U |
| 79-01-6 | Trichloroethylene | 1.55 | 7.3 | |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 1.55 | 1.6 | |
| 108-05-4 | Vinyl acetate | 1.55 | 0.54 | U |
| 593-60-2 | Vinyl bromide | 1.55 | 0.68 | U |
| 75-01-4 | Vinyl Chloride | 1.55 | 0.20 | U |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 219973 | 12.062 | 207186 | 12.072 | |
| ISTD: 1,4-Difluorobenzene | 1127831 | 13.638 | 1080437 | 13.641 | |
| ISTD: d5-Chlorobenzene | 1059994 | 18.895 | 1034699 | 18.895 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. SDG: 21D0660
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 21D0660-02 File ID: TQ214403.D
 Sampled: 04/13/21 15:00 Prepared: 04/20/21 12:32 Analyzed: 04/21/21 00:28
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BD11066 Sequence: Y1D2135 Calibration: YC10005 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.75 | 1.2 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.75 | 0.96 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.75 | 1.2 | U |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 1.75 | 1.3 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.75 | 0.96 | U |
| 75-34-3 | 1,1-Dichloroethane | 1.75 | 0.71 | U |
| 75-35-4 | 1,1-Dichloroethylene | 1.75 | 0.17 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.75 | 1.3 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 1.75 | 5.3 | |
| 106-93-4 | 1,2-Dibromoethane | 1.75 | 1.3 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.75 | 1.1 | U |
| 107-06-2 | 1,2-Dichloroethane | 1.75 | 0.71 | U |
| 78-87-5 | 1,2-Dichloropropane | 1.75 | 0.81 | U |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 1.75 | 1.2 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 1.75 | 2.0 | |
| 106-99-0 | 1,3-Butadiene | 1.75 | 1.4 | |
| 541-73-1 | 1,3-Dichlorobenzene | 1.75 | 1.1 | U |
| 142-28-9 | 1,3-Dichloropropane | 1.75 | 0.81 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.75 | 1.1 | U |
| 123-91-1 | 1,4-Dioxane | 1.75 | 1.3 | U |
| 78-93-3 | 2-Butanone | 1.75 | 20 | |
| 591-78-6 | 2-Hexanone | 1.75 | 1.4 | U |
| 107-05-1 | 3-Chloropropene | 1.75 | 2.7 | U |
| 108-10-1 | 4-Methyl-2-pentanone | 1.75 | 3.0 | |
| 67-64-1 | Acetone | 1.75 | 65 | |
| 107-13-1 | Acrylonitrile | 1.75 | 0.38 | U |
| 71-43-2 | Benzene | 1.75 | 26 | |
| 100-44-7 | Benzyl chloride | 1.75 | 0.91 | U |
| 75-27-4 | Bromodichloromethane | 1.75 | 1.2 | U |
| 75-25-2 | Bromoform | 1.75 | 1.8 | U |
| 74-83-9 | Bromomethane | 1.75 | 0.68 | U |
| 75-15-0 | Carbon disulfide | 1.75 | 67 | |
| 56-23-5 | Carbon tetrachloride | 1.75 | 0.28 | U |
| 108-90-7 | Chlorobenzene | 1.75 | 0.81 | U |
| 75-00-3 | Chloroethane | 1.75 | 0.46 | U |
| 67-66-3 | Chloroform | 1.75 | 4.0 | |
| 74-87-3 | Chloromethane | 1.75 | 0.36 | U |
| 156-59-2 | cis-1,2-Dichloroethylene | 1.75 | 0.42 | |
| 10061-01-5 | cis-1,3-Dichloropropylene | 1.75 | 0.80 | U |
| 110-82-7 | Cyclohexane | 1.75 | 3.1 | |

Laboratory: York Analytical Laboratories, Inc. SDG: 21D0660
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 21D0660-02 File ID: TQ214403.D
 Sampled: 04/13/21 15:00 Prepared: 04/20/21 12:32 Analyzed: 04/21/21 00:28
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BD11066 Sequence: Y1D2135 Calibration: YC10005 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 124-48-1 | Dibromochloromethane | 1.75 | 1.5 | U |
| 75-71-8 | Dichlorodifluoromethane | 1.75 | 2.4 | |
| 141-78-6 | Ethyl acetate | 1.75 | 1.3 | U |
| 100-41-4 | Ethyl Benzene | 1.75 | 3.3 | |
| 87-68-3 | Hexachlorobutadiene | 1.75 | 1.9 | U |
| 67-63-0 | Isopropanol | 1.75 | 1.4 | |
| 80-62-6 | Methyl Methacrylate | 1.75 | 0.72 | U |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 1.75 | 0.63 | U |
| 75-09-2 | Methylene chloride | 1.75 | 1.9 | J+ |
| 142-82-5 | n-Heptane | 1.75 | 11 | |
| 110-54-3 | n-Hexane | 1.75 | 11 | |
| 95-47-6 | o-Xylene | 1.75 | 5.1 | |
| 179601-23-1 | p- & m- Xylenes | 1.75 | 11 | |
| 622-96-8 | p-Ethyltoluene | 1.75 | 5.1 | |
| 115-07-1 | Propylene | 1.75 | 36 | |
| 100-42-5 | Styrene | 1.75 | 0.75 | U |
| 127-18-4 | Tetrachloroethylene | 1.75 | 80 | |
| 109-99-9 | Tetrahydrofuran | 1.75 | 1.0 | U |
| 108-88-3 | Toluene | 1.75 | 19 | |
| 156-60-5 | trans-1,2-Dichloroethylene | 1.75 | 0.70 | U |
| 10061-02-6 | trans-1,3-Dichloropropylene | 1.75 | 0.80 | U |
| 79-01-6 | Trichloroethylene | 1.75 | 2.9 | |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 1.75 | 1.3 | |
| 108-05-4 | Vinyl acetate | 1.75 | 0.62 | U |
| 593-60-2 | Vinyl bromide | 1.75 | 0.77 | U |
| 75-01-4 | Vinyl Chloride | 1.75 | 0.22 | U |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 201198 | 12.066 | 232457 | 12.065 | |
| ISTD: 1,4-Difluorobenzene | 1021693 | 13.638 | 1189631 | 13.635 | |
| ISTD: d5-Chlorobenzene | 965801 | 18.888 | 1133083 | 18.888 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. SDG: 21D0660
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 21D0660-03 File ID: TQ214409.D
 Sampled: 04/13/21 15:00 Prepared: 04/20/21 12:32 Analyzed: 04/21/21 06:34
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BD11066 Sequence: Y1D2135 Calibration: YC10005 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.7 | 1.2 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.7 | 0.93 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.7 | 1.2 | U |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 1.7 | 1.3 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.7 | 0.93 | U |
| 75-34-3 | 1,1-Dichloroethane | 1.7 | 0.69 | U |
| 75-35-4 | 1,1-Dichloroethylene | 1.7 | 0.17 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.7 | 1.3 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 1.7 | 3.0 | |
| 106-93-4 | 1,2-Dibromoethane | 1.7 | 1.3 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.7 | 1.0 | U |
| 107-06-2 | 1,2-Dichloroethane | 1.7 | 0.69 | U |
| 78-87-5 | 1,2-Dichloropropane | 1.7 | 0.79 | U |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 1.7 | 1.2 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 1.7 | 0.84 | U |
| 106-99-0 | 1,3-Butadiene | 1.7 | 1.1 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.7 | 1.0 | U |
| 142-28-9 | 1,3-Dichloropropane | 1.7 | 0.79 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.7 | 1.0 | U |
| 123-91-1 | 1,4-Dioxane | 1.7 | 1.2 | U |
| 78-93-3 | 2-Butanone | 1.7 | 1.4 | |
| 591-78-6 | 2-Hexanone | 1.7 | 1.4 | U |
| 107-05-1 | 3-Chloropropene | 1.7 | 2.7 | U |
| 108-10-1 | 4-Methyl-2-pentanone | 1.7 | 0.70 | U |
| 67-64-1 | Acetone | 1.7 | 9.9 | |
| 107-13-1 | Acrylonitrile | 1.7 | 0.37 | U |
| 71-43-2 | Benzene | 1.7 | 0.54 | U |
| 100-44-7 | Benzyl chloride | 1.7 | 0.88 | U |
| 75-27-4 | Bromodichloromethane | 1.7 | 4.1 | |
| 75-25-2 | Bromoform | 1.7 | 1.8 | U |
| 74-83-9 | Bromomethane | 1.7 | 0.66 | U |
| 75-15-0 | Carbon disulfide | 1.7 | 1.3 | |
| 56-23-5 | Carbon tetrachloride | 1.7 | 0.64 | |
| 108-90-7 | Chlorobenzene | 1.7 | 0.78 | U |
| 75-00-3 | Chloroethane | 1.7 | 0.45 | U |
| 67-66-3 | Chloroform | 1.7 | 67 | |
| 74-87-3 | Chloromethane | 1.7 | 0.35 | U |
| 156-59-2 | cis-1,2-Dichloroethylene | 1.7 | 0.17 | U |
| 10061-01-5 | cis-1,3-Dichloropropylene | 1.7 | 0.77 | U |
| 110-82-7 | Cyclohexane | 1.7 | 0.59 | U |

Laboratory: York Analytical Laboratories, Inc. SDG: 21D0660
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 21D0660-03 File ID: TQ214409.D
 Sampled: 04/13/21 15:00 Prepared: 04/20/21 12:32 Analyzed: 04/21/21 06:34
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BD11066 Sequence: Y1D2135 Calibration: YC10005 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|---|
| 124-48-1 | Dibromochloromethane | 1.7 | 1.4 | U |
| 75-71-8 | Dichlorodifluoromethane | 1.7 | 2.3 | |
| 141-78-6 | Ethyl acetate | 1.7 | 1.2 | U |
| 100-41-4 | Ethyl Benzene | 1.7 | 0.81 | |
| 87-68-3 | Hexachlorobutadiene | 1.7 | 1.8 | U |
| 67-63-0 | Isopropanol | 1.7 | 0.84 | U |
| 80-62-6 | Methyl Methacrylate | 1.7 | 0.70 | U |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 1.7 | 0.61 | U |
| 75-09-2 | Methylene chloride | 1.7 | 1.2 | U |
| 142-82-5 | n-Heptane | 1.7 | 0.70 | U |
| 110-54-3 | n-Hexane | 1.7 | 0.78 | |
| 95-47-6 | o-Xylene | 1.7 | 1.3 | |
| 179601-23-1 | p- & m- Xylenes | 1.7 | 3.9 | |
| 622-96-8 | p-Ethyltoluene | 1.7 | 2.0 | |
| 115-07-1 | Propylene | 1.7 | 1.8 | |
| 100-42-5 | Styrene | 1.7 | 0.73 | U |
| 127-18-4 | Tetrachloroethylene | 1.7 | 12 | |
| 109-99-9 | Tetrahydrofuran | 1.7 | 1.0 | U |
| 108-88-3 | Toluene | 1.7 | 1.6 | |
| 156-60-5 | trans-1,2-Dichloroethylene | 1.7 | 0.67 | U |
| 10061-02-6 | trans-1,3-Dichloropropylene | 1.7 | 0.77 | U |
| 79-01-6 | Trichloroethylene | 1.7 | 0.23 | U |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 1.7 | 1.3 | |
| 108-05-4 | Vinyl acetate | 1.7 | 0.60 | U |
| 593-60-2 | Vinyl bromide | 1.7 | 0.74 | U |
| 75-01-4 | Vinyl Chloride | 1.7 | 0.22 | U |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 242722 | 12.062 | 232457 | 12.065 | |
| ISTD: 1,4-Difluorobenzene | 1231707 | 13.638 | 1189631 | 13.635 | |
| ISTD: d5-Chlorobenzene | 1124826 | 18.892 | 1133083 | 18.888 | |

* Values outside of QC limits

JR 12/8/2025

December 8, 2025

Mr. Jason Stewart
Advanced Cleanup Technologies
228 Park Ave S PMB 34864
New York, New York 10003

Re: Data Usability Summary Report – York Analytical Laboratories, Inc. - 22F1004

Dear Mr. Stewart:

The evaluation of analytical data by York Analytical Laboratories for project 9268-BXNY, which were reported in a single data package under Job No. 22F1004 has been completed. The following samples were reported.

SS-15 OA-3 IA-14

Analysis was performed in accordance with EPA Method TO-15 (volatile organics). The review was performed to the extent possible, in accordance with the analytical method, and “DER-10/ Technical Guidance for Site Investigation and Remediation”. Professional judgment is applied as necessary and appropriate. National Functional Guidelines for Organic Data Review was consulted as needed. Qualifiers consistent with those defined by EPA Region 2 are applied as necessary and appropriate.

Below is the Data Usability Summary Report (DUSR) associated with these samples.

| Data Usability Summary Report | |
|---|----------------------------|
| 1. Is the data package complete as defined under the requirements for the most current DEC ASP Category B or USEPA CLP data deliverables? | Yes |
| 2. Have all holding times been met? | Yes |
| 3. Do all the QC data; blanks, instrument tunings, calibration standards, calibration verifications, surrogate recoveries, spike recoveries, replicate analyses, laboratory controls and sample data fall within the protocol required limits and specifications? | No -see following sections |
| 4. Have all of the data been generated using established and agreed upon analytical protocols? | Yes |
| 5. Does an evaluation of the raw data confirm the results provided in the data summary sheet and the quality control verification forms? | Yes |
| 6. Have the correct data qualifiers been used and are they consistent with the most current DEC ASP? | Yes |
| 7. Have any quality control (QC) exceedances been specifically noted in the DUSR and have the corresponding QC summary sheet from the data package been attached to the DUSR? | Yes |

Overall Evaluation

Based on the data review effort, results are usable, with the following qualifications. For samples that are qualified as estimated (J-, UJ), detected results may be biased low. False negatives may exist in non-detect results. Sample results that are qualified as estimated (J+) may be biased high. For samples that are qualified as estimated with any combination of (J), (J-) and/or (J+), the (J) qualifier takes precedence and is applied to the sample result. It is not possible to determine the direction of the bias and the overall effect on the result.

- The result for isopropanol in SS-15 and OA-13 is qualified as estimated (J-) due to unacceptable calibration.
- The results 1,2,4-trichlorobenzene, 2-hexanone, 4-methyl-2-pentanone and hexachlorobutadiene in all samples are qualified as estimated (UJ) due to low response in the continuing calibration verification (CCV) standard.
- The results for 1,2,4-trichlorobenzene, 2-hexanone, 4-methyl-2-pentanone and hexachlorobutadiene in all samples are qualified as estimated (UJ) due to recovery in the laboratory control sample (LCS).

Qualifier definitions are provided in Attachment A. A copy of the chain of custody record is provided in Attachment B. Pages from the data package illustrating the exceedances and issues described in this validation report are included in Attachment C. Annotated Summary Forms are included in Attachment D detailing qualifications resulting from the data review effort.

The following components were reviewed, where applicable:

- Chain of Custody
- Receiving conditions
- Holding times
- Preservation
- Analyte lists
- Reporting limits
- Requested methods
- Units, and
- Sample related quality control data:
 - Method, instrument blanks
 - Clean canister certification
 - Field blanks
 - Surrogate recoveries
 - LCS recoveries
 - Internal standard area response
 - Duplicates
 - Analyte Identification

- Instrument related quality control data:
 - Instrument tunes
 - Calibration summaries

The following sections of the report detail only quality control exceedances that impacted results. Where a quality control item exceeded control limits but there is no impact to the samples results, these are not detailed in the report.

Documentation: A completeness review of the data package was performed, and the data package was determined to be a complete Category B data package. The following documentation issues were observed during the review:

- The data package indicates that the laboratory did not hold certification for ethyl acetate, p-ethyltoluene, propylene, tetrahydrofuran, 1,1,1,2-tetrachloroethane, 1,3-dichloropropane, and 2-hexanone. It is noted that NY ELAP does not offer certification for ethyl acetate, p-ethyltoluene, propylene, tetrahydrofuran and 2-hexanone, but does offer certification for 1,1,1,2-tetrachloroethane and 1,3-dichloropropane. The laboratory was contacted to request clarification. The laboratory stated *'Given the age of the project, it is difficult to ascertain the exact circumstances in which these compounds were requested. We don't typically report non-certified compounds, except at the request of the client, so it can be safely assumed that there was correspondence regarding the accreditation status of the compounds at the time the samples were taken. Just to note- while these compounds are indeed non-certified under our scope of accreditation between 2020-2022, we ran the compounds in the exact same way we ran our certified compounds – quantitated against a multiple calibration, where they were properly qualified as specified by our certifying body.'*
- The Summary Form for the CCV analyzed on 06/22/22 06:51 incorrectly identifies the analysis time as 04/19/22 22:46.

Holding Times, Preservation, Sample Integrity:

A copy of the applicable chain of custody (COC) record was included in the data package, documenting sample collection dates of June 16 and 17, 2022. The samples were received at the laboratory on the same day on June 20, 2022. The samples were analyzed within method hold time. However, the data package does contain sample receipt information regarding condition of canisters upon receipt. The sample preparation bench sheets documents canister vacuum upon receipt, which indicates all canisters preserved (pressure inside the canister maintained within +/- 5 psi from sampling to check in the laboratory or analysis).

A. Volatile Organics

1. Calibration

One initial calibration (IC) was performed in support of the sample analysis. All relative

response factors (RRFs) and relative standard deviations (RSDs) or correlation coefficients are acceptable with the exception noted below.

| Analyte | RSD |
|-------------|------|
| Isopropanol | 55.2 |

The laboratory applied quadratic equations to achieve an acceptable correlation coefficient and used this model to quantitate sample concentrations. However, a review of the individual calibration levels indicates high responses in the lower calibration standards as presented below.

| Analyte | RRF 50 | RRF 30 | RRF 20 | RRF 10 | RRF 3 | RRF 0.5 | RRF 0.2 | RRF 0.1 |
|-------------|----------|----------|---------|----------|----------|-----------------|-----------------|-----------------|
| Isopropanol | 1.638682 | 1.676867 | 1.67942 | 1.781029 | 2.102878 | 2.991038 | 4.032331 | 5.756278 |

The high RSD for isopropanol is caused by excessively high response factors in the three lowest concentration standards (**bolded above**). The validator calculated the RSD with these two standards excluded, and the RSD for isopropanol is 10.7%.

The recalculated sample concentration using the updated average relative response factor is presented below:

| Sample | Reported Isopropanol (on-column ppbv) | Recalculated Isopropanol (on-column ppbv) | %D |
|--------|---------------------------------------|---|-----|
| SS-15 | 0.30 | 0.67 | 123 |
| OA-13 | 0.62 | 0.98 | 58 |
| IA-14 | 8.58 | 8.51 | 08 |

Where the difference in concentration is less than 5%, no action is taken. Where the difference in concentration exceeds 5% sample results are qualified as estimated (J-), biased low due to unacceptable calibration. The result for isopropanol in SS-15 and OA-13 is qualified as estimated (J-) due to unacceptable calibration. The concentration of isopropanol in IA-14 is greater than the concentration of the 3ppv standard and the high responses for isopropanol in the low standards has no impact on sample results.

A second source ICV standard was analyzed after the IC, and all percent differences are acceptable ($\leq 30\%D$). CCVs were analyzed at the appropriate frequency and are acceptable ($\%D \leq 30$) with the exceptions noted below.

| Analyte | %D | Associated Sample | Qualifier Applied |
|---------------------------|-------|-------------------|-------------------|
| <i>CCV 06/22/22 06:51</i> | | | |
| 1,2,4-Trichlorobenzene | -42.8 | OA-3 | UJ |
| 2-Hexanone | -22.4 | IA-14 | |
| 4-Methyl-2-pentanone | -42.0 | | |
| Hexachlorobutadiene | -41.8 | | |
| <i>CCV 04/19/22 22:46</i> | | | |
| 1,2,4-Trichlorobenzene | -36.1 | SS-15 | UJ |
| 2-Hexanone | -42.5 | | |
| 4-Methyl-2-pentanone | -39.6 | | |
| Hexachlorobutadiene | -39.3 | | |

The percent differences represent a decrease in instrument sensitivity. The results 1,2,4-trichlorobenzene, 2-hexanone, 4-methyl-2-pentanone and hexachlorobutadiene in all samples are qualified as estimated (UJ) due to low response in the CCV.

2. Laboratory Control Sample (LCS) /LCS Duplicate (LCSD)

Two LCSs were prepared and analyzed with the sample. All recoveries are acceptable (70-130%R), with the following exceptions:

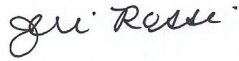
| Analyte | LCS %R | Affected Sample | Qualifier Applied |
|------------------------|--------|-----------------|-------------------|
| <i>BF21412-BS1</i> | | | |
| 1,2,4-Trichlorobenzene | 62.1 | OA-3 | UJ |
| 2-Hexanone | 53.4 | IA-14 | |
| 4-Methyl-2-pentanone | 56.6 | | |
| Hexachlorobutadiene | 57.2 | | |
| Analyte | LCS %R | Affected Sample | Qualifier Applied |
| <i>BF21498-BS1</i> | | | |
| 1,2,4-Trichlorobenzene | 68.2 | SS-15 | UJ |
| 2-Hexanone | 57.4 | | |
| 4-Methyl-2-pentanone | 60.0 | | |
| Hexachlorobutadiene | 62.3 | | |

The results for 1,2,4-trichlorobenzene, 2-hexanone, 4-methyl-2-pentanone and hexachlorobutadiene in all samples are qualified as estimated (UJ) due to recovery in the LCS.

22F1004

No other sample results are qualified. Please feel free to contact me at (908) 370-3431 or richjერიrossi513@gmail.com if you have any questions regarding this data package review report or need further information.

Sincerely,

A handwritten signature in cursive script that reads "Jeri Rossi". The signature is written in black ink on a white background.

Jeri L Rossi, CEAC

Environmental Consulting Chemist

ATTACHMENT A

Qualifier Definitions

EPA Qualifier Definitions

- U The analyte was analyzed for but was not detected above the level of the reported sample quantitation limit.
- J The result is an estimated quantity. The associated numerical value is the approximate concentration of the analyte in the sample.
- J+ The result is an estimated quantity, but the result may be biased high.
- J- The result is an estimated quantity, but the result may be biased low.
- NJ The analyte has been “tentatively identified” or “presumptively” as present and the associated numerical value is the estimated concentration in the sample.
- UJ The analyte was analyzed for but was not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
- R The data are unusable. The sample results are rejected due to serious deficiencies in meeting QC criteria. The analyte may or may not be present in the sample.

ATTACHMENT B

CHAIN OF CUSTODY (COC)



York Analytical Laboratories, Inc.
120 Research Drive Stratford, CT 06615
132-02 89th Ave Queens, NY 11418

Field Chain-of-Custody Record - AIR

YORK Project No.
22F1004



clientservices@yorklab.com
www.yorklab.com

NOTE: YORK's Standard Terms & Conditions are listed on the back side of this document. This document serves as your written authorization for YORK to proceed with the analyses requested below. signature binds you to YORK's Standard Terms & Conditions.

Your Page ___ of ___

| | | | | | | | | | |
|---|--|--------------------------------|--|--------------------------------|--|----------------------------|--|---|--|
| YOUR Information | | Report To: | | Invoice To: | | YOUR Project Number | | Turn-Around Time | |
| Company: Advanced Cleanup Tech | | Company: ACT | | Company: ACT | | 9628-BXNY | | RUSH - Next Day | |
| Address: 228 Park Ave S PMB 3484 | | Address: same | | Address: same | | YOUR Project Name | | RUSH - Two Day | |
| Address: NYNY 10003 | | | | | | | | RUSH - Three Day | |
| Phone: 631 466 4938 | | Phone: | | Phone: | | | | RUSH - Four Day | |
| Contact: Jason Stewart | | Contact: Paul Stewart | | Contact: Kera Friedman | | | | Standard (5-7 Day) <input checked="" type="checkbox"/> | |
| E-mail: jasons@act-earth | | E-mail: Pauls@act-earth | | E-mail: Keraf@act-earth | | YOUR PO#: | | | |

Please print clearly and legibly. All information must be complete. Samples will not be logged in and the turn-around-time clock will not begin until any questions by YORK are resolved.

| | | | | | | |
|---|--|--|--|----------------------|--|---|
| Samples Collected by: (print your name above and sign below) Tim Yang | Air Matrix Codes | Samples From | Report / EDD Type (circle selections) | | | YORK Reg. Comp. Compared to the following Regulation(s): (please fill in) |
| | <input checked="" type="radio"/> AI - Indoor Ambient Air | New York | <input checked="" type="checkbox"/> Summary Report | CT RCP | <input checked="" type="checkbox"/> Standard Excel EDD | |
| | <input checked="" type="radio"/> AO - Outdoor Amb. Air | New Jersey | QA Report | CT RCP DQA/DUE | EQuIS (Standard) | |
| | AE - Vapor Extraction Well/ Process Gas/Effluent | Connecticut | NY ASP A Package | NJDEP Reduced Deliv. | NYSDEC EQuIS | |
| | Pennsylvania | <input checked="" type="checkbox"/> NY ASP B Package | NJDQKP | NJDEP SRP HazSite | | |
| | Other | Other: | | | | |

Certified Canisters: Batch ___ Individual ___ **Please enter the following REQUIRED Field Data** Reporting Units: ug/m³ ___ ppbv ___ ppmv ___

| Sample Identification | Date/Time Sampled | Air Matrix | Canister Vacuum Before Sampling (in Hg) | Canister Vacuum After Sampling (in Hg) | Canister ID | Flow Cont. ID | Analysis Requested |
|-----------------------|-------------------|------------|---|--|-------------|---------------|--------------------|
| SS-15 | 6/16/22 9-1 | AS | -30 | -9 | 36988 | 7093 | TO-15 |
| OA-3 | 6/16-6/17 | AO | -30 | -10 | 10013 | 7364 | " |
| IA-14 | " " | AI | -30 | -10 | 42991 | 5606 | " |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |
| | | | | | | | |

| | | |
|------------------|--|--|
| Comments: | Detection Limits Required | Sampling Media |
| | ≤ 1 ug/m ³ <input checked="" type="checkbox"/> NYSDEC V1 Limits ___ Routine Survey ___ Other ___ | 6 Liter Canister <input checked="" type="checkbox"/> Tedlar Bag ___ |

| | | | | | |
|-----------------------------------|-----------------------|-----------------------------------|-----------------------|-----------------------------------|----------------------|
| Samples Relinquished by / Company | Date/Time | Samples Received by / Company | Date/Time | Samples Relinquished by / Company | Date/Time |
| ACT GARAGE | 6/20/22 7:10AM | KB Bunk York | 6/20/22 7:40AM | KB Bunk York | 6/20/22 16:52 |
| Samples Received by / Company | Date/Time | Samples Relinquished by / Company | Date/Time | Samples Received by / Company | Date/Time |
| 7 Gal | 6/20/22 16:52 | 7 Gal | 6/21/22 12:45 | Ivan B | 6-21-22 12:45 |
| Samples Relinquished by / Company | Date/Time | Samples Received in LAB by | Date/Time | | |
| Ivan B | 6-21-22 | Akene A Schuck | 6/21/22 16:15 | | |

33 of 304

ATTACHMENT C

**SELECTED PAGES FROM DATA PACKAGE –
QC EXCEEDANCES AND VALIDATION ISSUES**

FORM III

LCS / LCS DUPLICATE RECOVERY

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 22F1004Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYMatrix: AirBatch: BF21412Laboratory ID: BF21412-BS1Preparation: EPA TO15 PREPInitial/Final: 400 mL / 400 mL

| COMPOUND | SPIKE ADDED (ppbv) | LCS CONCENTRATION (ppbv) | LCS % REC. # | QC LIMITS REC. |
|---|--------------------|--------------------------|--------------|----------------|
| 1,1,1,2-Tetrachloroethane | 10.0 | 8.89 | 88.9 | 70 - 130 |
| 1,1,1-Trichloroethane | 10.0 | 10.3 | 103 | 70 - 130 |
| 1,1,2,2-Tetrachloroethane | 10.0 | 7.92 | 79.2 | 70 - 130 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 10.0 | 10.1 | 101 | 70 - 130 |
| 1,1,2-Trichloroethane | 10.0 | 8.39 | 83.9 | 70 - 130 |
| 1,1-Dichloroethane | 10.0 | 9.19 | 91.9 | 70 - 130 |
| 1,1-Dichloroethylene | 10.0 | 8.73 | 87.3 | 70 - 130 |
| 1,2,4-Trichlorobenzene | 10.0 | 6.21 | 62.1 * | 70 - 130 |
| 1,2,4-Trimethylbenzene | 10.0 | 7.83 | 78.3 | 70 - 130 |
| 1,2-Dibromoethane | 10.0 | 8.47 | 84.7 | 70 - 130 |
| 1,2-Dichlorobenzene | 10.0 | 7.51 | 75.1 | 70 - 130 |
| 1,2-Dichloroethane | 10.0 | 9.32 | 93.2 | 70 - 130 |
| 1,2-Dichloropropane | 10.0 | 7.59 | 75.9 | 70 - 130 |
| 1,2-Dichlorotetrafluoroethane | 10.0 | 11.2 | 112 | 70 - 130 |
| 1,3,5-Trimethylbenzene | 10.0 | 7.93 | 79.3 | 70 - 130 |
| 1,3-Butadiene | 10.0 | 7.28 | 72.8 | 70 - 130 |
| 1,3-Dichlorobenzene | 10.0 | 7.71 | 77.1 | 70 - 130 |
| 1,3-Dichloropropane | 10.0 | 8.06 | 80.6 | 70 - 130 |
| 1,4-Dichlorobenzene | 10.0 | 7.55 | 75.5 | 70 - 130 |
| 1,4-Dioxane | 10.0 | 7.04 | 70.4 | 70 - 130 |
| 2-Butanone | 10.0 | 7.77 | 77.7 | 70 - 130 |
| 2-Hexanone | 10.0 | 5.34 | 53.4 * | 70 - 130 |
| 3-Chloropropene | 10.0 | 7.83 | 78.3 | 70 - 130 |
| 4-Methyl-2-pentanone | 10.0 | 5.66 | 56.6 * | 70 - 130 |
| Acetone | 10.0 | 7.04 | 70.4 | 70 - 130 |
| Acrylonitrile | 10.0 | 7.99 | 79.9 | 70 - 130 |
| Benzene | 10.0 | 9.11 | 91.1 | 70 - 130 |
| Benzyl chloride | 10.0 | 7.39 | 73.9 | 70 - 130 |
| Bromodichloromethane | 10.0 | 8.29 | 82.9 | 70 - 130 |
| Bromoform | 10.0 | 8.93 | 89.3 | 70 - 130 |

FORM III

LCS / LCS DUPLICATE RECOVERY

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 22F1004Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYMatrix: AirBatch: BF21412Laboratory ID: BF21412-BS1Preparation: EPA TO15 PREPInitial/Final: 400 mL / 400 mL

| COMPOUND | SPIKE ADDED (ppbv) | LCS CONCENTRATION (ppbv) | LCS % REC. # | QC LIMITS REC. |
|--------------------------------|--------------------------|--------------------------------|--------------------|----------------------|
| Bromomethane | 10.0 | 10.2 | 102 | 70 - 130 |
| Carbon disulfide | 10.0 | 9.59 | 95.9 | 70 - 130 |
| Carbon tetrachloride | 10.0 | 9.51 | 95.1 | 70 - 130 |
| Chlorobenzene | 10.0 | 8.36 | 83.6 | 70 - 130 |
| Chloroethane | 10.0 | 9.11 | 91.1 | 70 - 130 |
| Chloroform | 10.0 | 9.90 | 99.0 | 70 - 130 |
| Chloromethane | 10.0 | 7.18 | 71.8 | 70 - 130 |
| cis-1,2-Dichloroethylene | 10.0 | 8.52 | 85.2 | 70 - 130 |
| cis-1,3-Dichloropropylene | 10.0 | 8.19 | 81.9 | 70 - 130 |
| Cyclohexane | 10.0 | 8.75 | 87.5 | 70 - 130 |
| Dibromochloromethane | 10.0 | 9.36 | 93.6 | 70 - 130 |
| Dichlorodifluoromethane | 10.0 | 10.0 | 100 | 70 - 130 |
| Ethyl acetate | 10.0 | 8.17 | 81.7 | 70 - 130 |
| Ethyl Benzene | 10.0 | 8.55 | 85.5 | 70 - 130 |
| Hexachlorobutadiene | 10.0 | 5.72 | 57.2 * | 70 - 130 |
| Isopropanol | 10.0 | 8.18 | 81.8 | 70 - 130 |
| Methyl Methacrylate | 10.0 | 7.69 | 76.9 | 70 - 130 |
| Methyl tert-butyl ether (MTBE) | 10.0 | 9.50 | 95.0 | 70 - 130 |
| Methylene chloride | 10.0 | 7.70 | 77.0 | 70 - 130 |
| n-Heptane | 10.0 | 8.15 | 81.5 | 70 - 130 |
| n-Hexane | 10.0 | 8.75 | 87.5 | 70 - 130 |
| o-Xylene | 10.0 | 8.49 | 84.9 | 70 - 130 |
| p- & m- Xylenes | 20.0 | 17.1 | 85.7 | 70 - 130 |
| p-Ethyltoluene | 10.0 | 8.14 | 81.4 | 70 - 130 |
| Propylene | 10.0 | 7.82 | 78.2 | 70 - 130 |
| Styrene | 10.0 | 8.09 | 80.9 | 70 - 130 |
| Tetrachloroethylene | 10.0 | 7.78 | 77.8 | 70 - 130 |
| Tetrahydrofuran | 10.0 | 7.83 | 78.3 | 70 - 130 |
| Toluene | 10.0 | 8.11 | 81.1 | 70 - 130 |
| trans-1,2-Dichloroethylene | 10.0 | 9.06 | 90.6 | 70 - 130 |

FORM III

LCS / LCS DUPLICATE RECOVERY

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 22F1004Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYMatrix: AirBatch: BF21498Laboratory ID: BF21498-BS1Preparation: EPA TO15 PREPInitial/Final: 400 mL / 400 mL

| COMPOUND | SPIKE ADDED (ppbv) | LCS CONCENTRATION (ppbv) | LCS % REC. # | QC LIMITS REC. |
|---|--------------------|--------------------------|--------------|----------------|
| 1,1,1,2-Tetrachloroethane | 10.0 | 9.55 | 95.5 | 70 - 130 |
| 1,1,1-Trichloroethane | 10.0 | 11.4 | 114 | 70 - 130 |
| 1,1,2,2-Tetrachloroethane | 10.0 | 8.62 | 86.2 | 70 - 130 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 10.0 | 11.2 | 112 | 70 - 130 |
| 1,1,2-Trichloroethane | 10.0 | 9.14 | 91.4 | 70 - 130 |
| 1,1-Dichloroethane | 10.0 | 10.0 | 100 | 70 - 130 |
| 1,1-Dichloroethylene | 10.0 | 9.65 | 96.5 | 70 - 130 |
| 1,2,4-Trichlorobenzene | 10.0 | 6.82 | 68.2 * | 70 - 130 |
| 1,2,4-Trimethylbenzene | 10.0 | 8.72 | 87.2 | 70 - 130 |
| 1,2-Dibromoethane | 10.0 | 9.18 | 91.8 | 70 - 130 |
| 1,2-Dichlorobenzene | 10.0 | 8.29 | 82.9 | 70 - 130 |
| 1,2-Dichloroethane | 10.0 | 10.5 | 105 | 70 - 130 |
| 1,2-Dichloropropane | 10.0 | 7.98 | 79.8 | 70 - 130 |
| 1,2-Dichlorotetrafluoroethane | 10.0 | 12.4 | 124 | 70 - 130 |
| 1,3,5-Trimethylbenzene | 10.0 | 8.84 | 88.4 | 70 - 130 |
| 1,3-Butadiene | 10.0 | 8.10 | 81.0 | 70 - 130 |
| 1,3-Dichlorobenzene | 10.0 | 8.54 | 85.4 | 70 - 130 |
| 1,3-Dichloropropane | 10.0 | 8.66 | 86.6 | 70 - 130 |
| 1,4-Dichlorobenzene | 10.0 | 8.31 | 83.1 | 70 - 130 |
| 1,4-Dioxane | 10.0 | 7.51 | 75.1 | 70 - 130 |
| 2-Butanone | 10.0 | 8.37 | 83.7 | 70 - 130 |
| 2-Hexanone | 10.0 | 5.74 | 57.4 * | 70 - 130 |
| 3-Chloropropene | 10.0 | 8.24 | 82.4 | 70 - 130 |
| 4-Methyl-2-pentanone | 10.0 | 6.00 | 60.0 * | 70 - 130 |
| Acetone | 10.0 | 7.76 | 77.6 | 70 - 130 |
| Acrylonitrile | 10.0 | 8.48 | 84.8 | 70 - 130 |
| Benzene | 10.0 | 9.87 | 98.7 | 70 - 130 |
| Benzyl chloride | 10.0 | 8.32 | 83.2 | 70 - 130 |
| Bromodichloromethane | 10.0 | 9.01 | 90.1 | 70 - 130 |
| Bromoform | 10.0 | 9.87 | 98.7 | 70 - 130 |

FORM III

LCS / LCS DUPLICATE RECOVERY

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 22F1004Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYMatrix: AirBatch: BF21498Laboratory ID: BF21498-BS1Preparation: EPA TO15 PREPInitial/Final: 400 mL / 400 mL

| COMPOUND | SPIKE ADDED (ppbv) | LCS CONCENTRATION (ppbv) | LCS % REC. # | QC LIMITS REC. |
|--------------------------------|--------------------|--------------------------|--------------|----------------|
| Bromomethane | 10.0 | 11.4 | 114 | 70 - 130 |
| Carbon disulfide | 10.0 | 10.4 | 104 | 70 - 130 |
| Carbon tetrachloride | 10.0 | 10.0 | 100 | 70 - 130 |
| Chlorobenzene | 10.0 | 9.04 | 90.4 | 70 - 130 |
| Chloroethane | 10.0 | 9.93 | 99.3 | 70 - 130 |
| Chloroform | 10.0 | 11.0 | 110 | 70 - 130 |
| Chloromethane | 10.0 | 8.27 | 82.7 | 70 - 130 |
| cis-1,2-Dichloroethylene | 10.0 | 9.15 | 91.5 | 70 - 130 |
| cis-1,3-Dichloropropylene | 10.0 | 8.75 | 87.5 | 70 - 130 |
| Cyclohexane | 10.0 | 9.42 | 94.2 | 70 - 130 |
| Dibromochloromethane | 10.0 | 10.2 | 102 | 70 - 130 |
| Dichlorodifluoromethane | 10.0 | 11.2 | 112 | 70 - 130 |
| Ethyl acetate | 10.0 | 8.77 | 87.7 | 70 - 130 |
| Ethyl Benzene | 10.0 | 9.19 | 91.9 | 70 - 130 |
| Hexachlorobutadiene | 10.0 | 6.23 | 62.3 * | 70 - 130 |
| Isopropanol | 10.0 | 8.93 | 89.3 | 70 - 130 |
| Methyl Methacrylate | 10.0 | 8.27 | 82.7 | 70 - 130 |
| Methyl tert-butyl ether (MTBE) | 10.0 | 10.3 | 103 | 70 - 130 |
| Methylene chloride | 10.0 | 8.30 | 83.0 | 70 - 130 |
| n-Heptane | 10.0 | 8.58 | 85.8 | 70 - 130 |
| n-Hexane | 10.0 | 9.32 | 93.2 | 70 - 130 |
| o-Xylene | 10.0 | 9.26 | 92.6 | 70 - 130 |
| p- & m- Xylenes | 20.0 | 18.7 | 93.5 | 70 - 130 |
| p-Ethyltoluene | 10.0 | 9.15 | 91.5 | 70 - 130 |
| Propylene | 10.0 | 7.94 | 79.4 | 70 - 130 |
| Styrene | 10.0 | 8.84 | 88.4 | 70 - 130 |
| Tetrachloroethylene | 10.0 | 8.45 | 84.5 | 70 - 130 |
| Tetrahydrofuran | 10.0 | 8.35 | 83.5 | 70 - 130 |
| Toluene | 10.0 | 8.77 | 87.7 | 70 - 130 |
| trans-1,2-Dichloroethylene | 10.0 | 9.88 | 98.8 | 70 - 130 |

FORM VI

INITIAL CALIBRATION DATA (Continued)

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 22F1004Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYCalibration: YD20018Instrument: TO15_AIR2Calibration Date: 04/19/22 22:46

| Compound | Mean RF | RF RSD | Mean RT | RT RSD | Linear r | Quad COD | LIMIT | Q |
|--------------------------------|-----------|----------|----------|--------------|----------|-----------|-------|---|
| Carbon disulfide | 3.255539 | 6.903854 | 9.33625 | 6.506453E-02 | | | 30 | |
| Carbon tetrachloride | 2.382535 | 8.868214 | 12.94933 | 5.208245E-02 | | | 30 | |
| Chlorobenzene | 1.052086 | 7.010271 | 18.87275 | 2.334866E-02 | | | 30 | |
| Chloroethane | 0.506012 | 5.264119 | 6.77975 | 0.1984605 | | | 30 | |
| Chloroform | 2.668734 | 8.02208 | 11.74725 | 7.008658E-02 | | | 30 | |
| Chloromethane | 0.5259801 | 13.78232 | 5.45325 | 0.2028036 | | | 30 | |
| cis-1,2-Dichloroethylene | 1.576247 | 12.11326 | 11.4747 | 7.087928E-02 | | | 30 | |
| cis-1,3-Dichloropropylene | 0.4575901 | 8.479898 | 15.6665 | 1.989716E-02 | | | 30 | |
| Cyclohexane | 1.625653 | 7.465707 | 12.65025 | 5.826469E-02 | | | 30 | |
| Dibromochloromethane | 0.5761676 | 15.29481 | 17.70325 | 1.921471E-02 | | | 30 | |
| Dichlorodifluoromethane | 3.200946 | 10.39896 | 4.96725 | 0.1882287 | | | 30 | |
| Ethanol | | | | | | | 30 | |
| Ethyl acetate | 1.994603 | 6.988111 | 11.39713 | 4.629581E-02 | | | 30 | |
| Ethyl Benzene | 1.603083 | 8.647667 | 18.97925 | 0.0255301 | | | 30 | |
| Hexachlorobutadiene | 0.6539784 | 15.252 | 27.20943 | 8.072138E-03 | | | 30 | |
| Isopropanol | 2.707315 | 55.16935 | 7.882875 | 0.1645517 | | 0.999882 | 0.99 | |
| Isopropylbenzene | 1.793419 | 8.106145 | 20.55987 | 1.771929E-02 | | | 30 | |
| Methyl Methacrylate | 0.2877017 | 8.989548 | 14.43062 | 3.711034E-02 | | | 30 | |
| Methyl tert-butyl ether (MTBE) | 3.477153 | 7.260362 | 9.621125 | 6.637063E-02 | | | 30 | |
| Methylene chloride | 1.130144 | 21.85847 | 9.20875 | 0.1207584 | | | 30 | |
| Naphthalene | 2.302281 | 32.40631 | 27.324 | 1.406847E-02 | | 0.9995157 | 0.99 | |
| n-Butylbenzene | 1.631939 | 9.284772 | 23.93113 | 1.055984E-02 | | | 30 | |
| n-Heptane | 1.595517 | 10.80162 | 13.17475 | 6.448263E-02 | | | 30 | |
| n-Hexane | 1.565854 | 7.02213 | 10.10287 | 7.638676E-02 | | | 30 | |
| n-Propylbenzene | 2.046093 | 7.358667 | 21.31875 | 2.267728E-02 | | | 30 | |
| o-Xylene | 1.265449 | 7.717153 | 19.9045 | 0.0200949 | | | 30 | |
| p- & m- Xylenes | 1.203784 | 11.61558 | 19.125 | 1.820726E-02 | | | 30 | |
| p-Ethyltoluene | 1.769097 | 5.460827 | 21.54 | 2.299606E-02 | | | 30 | |
| p-Isopropyltoluene | 1.840235 | 6.464565 | 23.12188 | 2.365521E-02 | | | 30 | |
| Propylene | 0.4830245 | 14.54994 | 4.893375 | 0.2201734 | | | 30 | |
| sec-Butylbenzene | 2.187856 | 7.075805 | 22.85425 | 2.336683E-02 | | | 30 | |
| Styrene | 1.018308 | 5.814062 | 19.92475 | 1.685623E-02 | | | 30 | |

FORM VII

CONTINUING CALIBRATION CHECK

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 22F1004Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYInstrument ID: TO15_AIR2Calibration: YD20018Lab File ID: TQ220717.DCalibration Date: 04/19/22 22:46Sequence: Y2F2411Injection Date: 06/22/22Lab Sample ID: Y2F2411-CCV1Injection Time: 06:51

| COMPOUND | TYPE | CONC. (ppbv) | | RESPONSE FACTOR | | | % DIFF / DRIFT | |
|---|------|--------------|------|-----------------|-----------|---------|----------------|-----------|
| | | STD | CCV | ICAL | CCV | MIN (#) | CCV | LIMIT (#) |
| 1,1,1,2-Tetrachloroethane | A | 10.0 | 8.76 | 0.5225608 | 0.4575546 | | -12.4 | 30 |
| 1,1,1-Trichloroethane | A | 10.0 | 10.2 | 2.761948 | 2.810861 | | 1.8 | 30 |
| 1,1,2,2-Tetrachloroethane | A | 10.0 | 8.04 | 0.8626136 | 0.6931939 | | -19.6 | 30 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | A | 10.0 | 10.1 | 2.475865 | 2.506197 | | 1.2 | 30 |
| 1,1,2-Trichloroethane | A | 10.0 | 8.43 | 0.3641926 | 0.3070471 | | -15.7 | 30 |
| 1,1-Dichloroethane | A | 10.0 | 9.10 | 2.022685 | 1.840634 | | -9.0 | 30 |
| 1,1-Dichloroethylene | A | 10.0 | 8.65 | 1.828947 | 1.581877 | | -13.5 | 30 |
| 1,2,4-Trichlorobenzene | A | 10.0 | 5.72 | 0.6607203 | 0.378042 | | -42.8 | 30 * |
| 1,2,4-Trimethylbenzene | A | 10.0 | 8.16 | 1.52027 | 1.24006 | | -18.4 | 30 |
| 1,2-Dibromoethane | A | 10.0 | 8.53 | 0.5731624 | 0.4887991 | | -14.7 | 30 |
| 1,2-Dichlorobenzene | A | 10.0 | 7.76 | 0.9529005 | 0.7392303 | | -22.4 | 30 |
| 1,2-Dichloroethane | A | 10.0 | 9.29 | 1.647525 | 1.5308 | | -7.1 | 30 |
| 1,2-Dichloropropane | A | 10.0 | 7.63 | 0.2644825 | 0.201719 | | -23.7 | 30 |
| 1,2-Dichlorotetrafluoroethane | A | 10.0 | 11.1 | 2.655057 | 2.938376 | | 10.7 | 30 |
| 1,3,5-Trimethylbenzene | A | 10.0 | 8.18 | 1.491271 | 1.220093 | | -18.2 | 30 |
| 1,3-Butadiene | A | 10.0 | 7.24 | 0.4907221 | 0.3551287 | | -27.6 | 30 |
| 1,3-Dichlorobenzene | A | 10.0 | 7.97 | 0.9521438 | 0.7585303 | | -20.3 | 30 |
| 1,3-Dichloropropane | A | 10.0 | 8.06 | 0.500202 | 0.4030905 | | -19.4 | 30 |
| 1,4-Dichlorobenzene | A | 10.0 | 7.87 | 0.9541104 | 0.7512614 | | -21.3 | 30 |
| 1,4-Dioxane | A | 10.0 | 7.30 | 0.2366858 | 0.1727497 | | -27.0 | 30 |
| 2-Butanone | A | 10.0 | 7.76 | 1.877054 | 1.45608 | | -22.4 | 30 |
| 2-Hexanone | A | 10.0 | 5.62 | 0.5317352 | 0.2988719 | | -43.8 | 30 * |
| 3-Chloropropene | A | 10.0 | 7.77 | 1.090371 | 0.8474979 | | -22.3 | 30 |
| 4-Methyl-2-pentanone | A | 10.0 | 5.80 | 0.5611524 | 0.325421 | | -42.0 | 30 * |
| Acetone | A | 10.0 | 7.02 | 1.624941 | 1.141238 | | -29.8 | 30 |
| Acrylonitrile | A | 10.0 | 7.72 | 0.7267274 | 0.5606935 | | -22.8 | 30 |
| Benzene | A | 10.0 | 9.12 | 3.988265 | 3.636066 | | -8.8 | 30 |
| Benzyl chloride | A | 10.0 | 7.56 | 1.118416 | 0.8453578 | | -24.4 | 30 |
| Bromodichloromethane | A | 10.0 | 8.36 | 0.5662355 | 0.4731749 | | -16.4 | 30 |

FORM VII

CONTINUING CALIBRATION CHECK

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 22F1004Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYInstrument ID: TO15_AIR2Calibration: YD20018Lab File ID: TQ220717.DCalibration Date: 04/19/22 22:46Sequence: Y2F2411Injection Date: 06/22/22Lab Sample ID: Y2F2411-CCV1Injection Time: 06:51

| COMPOUND | TYPE | CONC. (ppbv) | | RESPONSE FACTOR | | | % DIFF / DRIFT | |
|--------------------------------|------|--------------|------|-----------------|-----------|---------|----------------|-----------|
| | | STD | CCV | ICAL | CCV | MIN (#) | CCV | LIMIT (#) |
| Bromoform | A | 10.0 | 9.08 | 0.5589733 | 0.5073656 | | -9.2 | 30 |
| Bromomethane | A | 10.0 | 10.2 | 1.07543 | 1.094078 | | 1.7 | 30 |
| Carbon disulfide | A | 10.0 | 9.49 | 3.255539 | 3.089899 | | -5.1 | 30 |
| Carbon tetrachloride | A | 10.0 | 9.51 | 2.382535 | 2.301725 | | -3.4 | 30 |
| Chlorobenzene | A | 10.0 | 8.36 | 1.052086 | 0.880038 | | -16.4 | 30 |
| Chloroethane | A | 10.0 | 9.02 | 0.506012 | 0.4564789 | | -9.8 | 30 |
| Chloroform | A | 10.0 | 9.87 | 2.668734 | 2.634392 | | -1.3 | 30 |
| Chloromethane | A | 10.0 | 7.32 | 0.5259801 | 0.385212 | | -26.8 | 30 |
| cis-1,2-Dichloroethylene | A | 10.0 | 8.64 | 1.576247 | 1.361592 | | -13.6 | 30 |
| cis-1,3-Dichloropropylene | A | 10.0 | 8.16 | 0.4575901 | 0.3733286 | | -18.4 | 30 |
| Cyclohexane | A | 10.0 | 8.81 | 1.625653 | 1.432123 | | -11.9 | 30 |
| Dibromochloromethane | A | 10.0 | 9.38 | 0.5761676 | 0.5407028 | | -6.2 | 30 |
| Dichlorodifluoromethane | A | 10.0 | 10.0 | 3.200946 | 3.212941 | | 0.4 | 30 |
| Ethyl acetate | A | 10.0 | 8.18 | 1.994603 | 1.631095 | | -18.2 | 30 |
| Ethyl Benzene | A | 10.0 | 8.50 | 1.603083 | 1.36254 | | -15.0 | 30 |
| Hexachlorobutadiene | A | 10.0 | 5.99 | 0.6539784 | 0.3809212 | | -41.8 | 30 * |
| Isopropanol | Q | 10.0 | 8.41 | 2.707315 | 1.482894 | | -15.9 | 30 |
| Methyl Methacrylate | A | 10.0 | 7.75 | 0.2877017 | 0.2228967 | | -22.5 | 30 |
| Methyl tert-butyl ether (MTBE) | A | 10.0 | 9.47 | 3.477153 | 3.2913 | | -5.3 | 30 |
| Methylene chloride | A | 10.0 | 7.65 | 1.130144 | 0.8651118 | | -23.5 | 30 |
| n-Heptane | A | 10.0 | 8.23 | 1.595517 | 1.313021 | | -17.7 | 30 |
| n-Hexane | A | 10.0 | 8.77 | 1.565854 | 1.373648 | | -12.3 | 30 |
| o-Xylene | A | 10.0 | 8.47 | 1.265449 | 1.071284 | | -15.3 | 30 |
| p- & m- Xylenes | A | 20.0 | 17.1 | 1.203784 | 1.031113 | | -14.3 | 30 |
| p-Ethyltoluene | A | 10.0 | 8.39 | 1.769097 | 1.483905 | | -16.1 | 30 |
| Propylene | A | 10.0 | 7.77 | 0.4830245 | 0.3755248 | | -22.3 | 30 |
| Styrene | A | 10.0 | 8.17 | 1.018308 | 0.8319181 | | -18.3 | 30 |
| Tetrachloroethylene | A | 10.0 | 7.93 | 0.5641022 | 0.4474602 | | -20.7 | 30 |
| Tetrahydrofuran | A | 10.0 | 7.96 | 1.030652 | 0.8201629 | | -20.4 | 30 |

FORM VII

CONTINUING CALIBRATION CHECK

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 22F1004Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYInstrument ID: TO15_AIR2Calibration: YD20018Lab File ID: TQ220732.DCalibration Date: 04/19/22 22:46Sequence: Y2F2413Injection Date: 06/23/22Lab Sample ID: Y2F2413-CCV1Injection Time: 03:41

| COMPOUND | TYPE | CONC. (ppbv) | | RESPONSE FACTOR | | | % DIFF / DRIFT | |
|---|------|--------------|------|-----------------|-----------|---------|----------------|-----------|
| | | STD | CCV | ICAL | CCV | MIN (#) | CCV | LIMIT (#) |
| 1,1,1,2-Tetrachloroethane | A | 10.0 | 9.72 | 0.5225608 | 0.5079412 | | -2.8 | 30 |
| 1,1,1-Trichloroethane | A | 10.0 | 11.7 | 2.761948 | 3.2286 | | 16.9 | 30 |
| 1,1,2,2-Tetrachloroethane | A | 10.0 | 8.65 | 0.8626136 | 0.7462734 | | -13.5 | 30 |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | A | 10.0 | 11.7 | 2.475865 | 2.890816 | | 16.8 | 30 |
| 1,1,2-Trichloroethane | A | 10.0 | 9.17 | 0.3641926 | 0.3340704 | | -8.3 | 30 |
| 1,1-Dichloroethane | A | 10.0 | 10.2 | 2.022685 | 2.062586 | | 2.0 | 30 |
| 1,1-Dichloroethylene | A | 10.0 | 9.81 | 1.828947 | 1.794147 | | -1.9 | 30 |
| 1,2,4-Trichlorobenzene | A | 10.0 | 6.39 | 0.6607203 | 0.4221041 | | -36.1 | 30 * |
| 1,2,4-Trimethylbenzene | A | 10.0 | 8.69 | 1.52027 | 1.321068 | | -13.1 | 30 |
| 1,2-Dibromoethane | A | 10.0 | 9.24 | 0.5731624 | 0.5298719 | | -7.6 | 30 |
| 1,2-Dichlorobenzene | A | 10.0 | 8.17 | 0.9529005 | 0.7782002 | | -18.3 | 30 |
| 1,2-Dichloroethane | A | 10.0 | 10.8 | 1.647525 | 1.772224 | | 7.6 | 30 |
| 1,2-Dichloropropane | A | 10.0 | 8.09 | 0.2644825 | 0.2139312 | | -19.1 | 30 |
| 1,2-Dichlorotetrafluoroethane | A | 10.0 | 11.4 | 2.655057 | 3.022934 | | 13.9 | 30 |
| 1,3,5-Trimethylbenzene | A | 10.0 | 8.99 | 1.491271 | 1.340846 | | -10.1 | 30 |
| 1,3-Butadiene | A | 10.0 | 8.23 | 0.4907221 | 0.4040612 | | -17.7 | 30 |
| 1,3-Dichlorobenzene | A | 10.0 | 8.33 | 0.9521438 | 0.7929265 | | -16.7 | 30 |
| 1,3-Dichloropropane | A | 10.0 | 8.75 | 0.500202 | 0.4377096 | | -12.5 | 30 |
| 1,4-Dichlorobenzene | A | 10.0 | 8.21 | 0.9541104 | 0.7832056 | | -17.9 | 30 |
| 1,4-Dioxane | A | 10.0 | 7.58 | 0.2366858 | 0.1793665 | | -24.2 | 30 |
| 2-Butanone | A | 10.0 | 8.50 | 1.877054 | 1.595461 | | -15.0 | 30 |
| 2-Hexanone | A | 10.0 | 5.75 | 0.5317352 | 0.3058063 | | -42.5 | 30 * |
| 3-Chloropropene | A | 10.0 | 8.41 | 1.090371 | 0.9165837 | | -15.9 | 30 |
| 4-Methyl-2-pentanone | A | 10.0 | 6.04 | 0.5611524 | 0.338866 | | -39.6 | 30 * |
| Acetone | A | 10.0 | 8.09 | 1.624941 | 1.313952 | | -19.1 | 30 |
| Acrylonitrile | A | 10.0 | 8.60 | 0.7267274 | 0.6249856 | | -14.0 | 30 |
| Benzene | A | 10.0 | 10.0 | 3.988265 | 4.007355 | | 0.5 | 30 |
| Benzyl chloride | A | 10.0 | 8.09 | 1.118416 | 0.9049434 | | -19.1 | 30 |
| Bromodichloromethane | A | 10.0 | 9.25 | 0.5662355 | 0.5234904 | | -7.5 | 30 |

FORM VII

CONTINUING CALIBRATION CHECK

EPA TO-15

Laboratory: York Analytical Laboratories, Inc.SDG: 22F1004Client: Advanced Cleanup Technologies, Inc.Project: 9628-BXNYInstrument ID: TO15_AIR2Calibration: YD20018Lab File ID: TQ220732.DCalibration Date: 04/19/22 22:46Sequence: Y2F2413Injection Date: 06/23/22Lab Sample ID: Y2F2413-CCV1Injection Time: 03:41

| COMPOUND | TYPE | CONC. (ppbv) | | RESPONSE FACTOR | | | % DIFF / DRIFT | |
|--------------------------------|------|--------------|------|-----------------|-----------|---------|----------------|-----------|
| | | STD | CCV | ICAL | CCV | MIN (#) | CCV | LIMIT (#) |
| Bromoform | A | 10.0 | 9.91 | 0.5589733 | 0.55404 | | -0.9 | 30 |
| Bromomethane | A | 10.0 | 11.8 | 1.07543 | 1.267357 | | 17.8 | 30 |
| Carbon disulfide | A | 10.0 | 10.6 | 3.255539 | 3.468478 | | 6.5 | 30 |
| Carbon tetrachloride | A | 10.0 | 10.4 | 2.382535 | 2.52275 | | 5.9 | 30 |
| Chlorobenzene | A | 10.0 | 9.18 | 1.052086 | 0.9659292 | | -8.2 | 30 |
| Chloroethane | A | 10.0 | 10.4 | 0.506012 | 0.523877 | | 3.5 | 30 |
| Chloroform | A | 10.0 | 11.2 | 2.668734 | 2.998605 | | 12.4 | 30 |
| Chloromethane | A | 10.0 | 8.22 | 0.5259801 | 0.4325853 | | -17.8 | 30 |
| cis-1,2-Dichloroethylene | A | 10.0 | 9.60 | 1.576247 | 1.513359 | | -4.0 | 30 |
| cis-1,3-Dichloropropylene | A | 10.0 | 8.93 | 0.4575901 | 0.4084585 | | -10.7 | 30 |
| Cyclohexane | A | 10.0 | 9.51 | 1.625653 | 1.546279 | | -4.9 | 30 |
| Dibromochloromethane | A | 10.0 | 10.4 | 0.5761676 | 0.5974072 | | 3.7 | 30 |
| Dichlorodifluoromethane | A | 10.0 | 11.5 | 3.200946 | 3.676141 | | 14.8 | 30 |
| Ethyl acetate | A | 10.0 | 8.97 | 1.994603 | 1.789026 | | -10.3 | 30 |
| Ethyl Benzene | A | 10.0 | 9.31 | 1.603083 | 1.491957 | | -6.9 | 30 |
| Hexachlorobutadiene | A | 10.0 | 6.24 | 0.6539784 | 0.3970507 | | -39.3 | 30 * |
| Isopropanol | Q | 10.0 | 9.12 | 2.707315 | 1.60133 | | -8.8 | 30 |
| Methyl Methacrylate | A | 10.0 | 8.32 | 0.2877017 | 0.2393558 | | -16.8 | 30 |
| Methyl tert-butyl ether (MTBE) | A | 10.0 | 10.6 | 3.477153 | 3.675739 | | 5.7 | 30 |
| Methylene chloride | A | 10.0 | 8.49 | 1.130144 | 0.9598998 | | -15.1 | 30 |
| n-Heptane | A | 10.0 | 8.72 | 1.595517 | 1.391757 | | -12.8 | 30 |
| n-Hexane | A | 10.0 | 9.58 | 1.565854 | 1.500119 | | -4.2 | 30 |
| o-Xylene | A | 10.0 | 9.38 | 1.265449 | 1.18645 | | -6.2 | 30 |
| p- & m- Xylenes | A | 20.0 | 18.9 | 1.203784 | 1.136696 | | -5.6 | 30 |
| p-Ethyltoluene | A | 10.0 | 9.03 | 1.769097 | 1.597231 | | -9.7 | 30 |
| Propylene | A | 10.0 | 7.97 | 0.4830245 | 0.3849322 | | -20.3 | 30 |
| Styrene | A | 10.0 | 8.89 | 1.018308 | 0.9051465 | | -11.1 | 30 |
| Tetrachloroethylene | A | 10.0 | 8.51 | 0.5641022 | 0.4803092 | | -14.9 | 30 |
| Tetrahydrofuran | A | 10.0 | 8.53 | 1.030652 | 0.8793586 | | -14.7 | 30 |

ATTACHMENT D

ANNOTATED SUMMARY FORMS

Laboratory: York Analytical Laboratories, Inc. SDG: 22F1004
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 22F1004-01 File ID: TQ220742.D
 Sampled: 06/16/22 13:00 Prepared: 06/23/22 02:00 Analyzed: 06/23/22 14:19
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BF21498 Sequence: Y2F2413 Calibration: YD20018 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|----|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.93 | 1.3 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.93 | 1.1 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.93 | 1.3 | U |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 1.93 | 1.5 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.93 | 1.1 | U |
| 75-34-3 | 1,1-Dichloroethane | 1.93 | 0.78 | U |
| 75-35-4 | 1,1-Dichloroethylene | 1.93 | 0.19 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.93 | 1.4 | UJ |
| 95-63-6 | 1,2,4-Trimethylbenzene | 1.93 | 45 | |
| 106-93-4 | 1,2-Dibromoethane | 1.93 | 1.5 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.93 | 1.2 | U |
| 107-06-2 | 1,2-Dichloroethane | 1.93 | 0.78 | U |
| 78-87-5 | 1,2-Dichloropropane | 1.93 | 0.89 | U |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 1.93 | 1.4 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 1.93 | 57 | |
| 106-99-0 | 1,3-Butadiene | 1.93 | 1.3 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.93 | 1.2 | U |
| 142-28-9 | 1,3-Dichloropropane | 1.93 | 0.89 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.93 | 1.2 | U |
| 123-91-1 | 1,4-Dioxane | 1.93 | 1.4 | U |
| 78-93-3 | 2-Butanone | 1.93 | 3.6 | |
| 591-78-6 | 2-Hexanone | 1.93 | 1.6 | UJ |
| 107-05-1 | 3-Chloropropene | 1.93 | 3.0 | U |
| 108-10-1 | 4-Methyl-2-pentanone | 1.93 | 0.79 | UJ |
| 67-64-1 | Acetone | 1.93 | 23 | |
| 107-13-1 | Acrylonitrile | 1.93 | 0.42 | U |
| 71-43-2 | Benzene | 1.93 | 2.4 | |
| 100-44-7 | Benzyl chloride | 1.93 | 1.0 | U |
| 75-27-4 | Bromodichloromethane | 1.93 | 1.3 | U |
| 75-25-2 | Bromoform | 1.93 | 2.0 | U |
| 74-83-9 | Bromomethane | 1.93 | 0.75 | U |
| 75-15-0 | Carbon disulfide | 1.93 | 0.60 | U |
| 56-23-5 | Carbon tetrachloride | 1.93 | 0.36 | |
| 108-90-7 | Chlorobenzene | 1.93 | 0.89 | U |
| 75-00-3 | Chloroethane | 1.93 | 0.51 | U |
| 67-66-3 | Chloroform | 1.93 | 9.2 | |
| 74-87-3 | Chloromethane | 1.93 | 0.40 | U |
| 156-59-2 | cis-1,2-Dichloroethylene | 1.93 | 0.19 | U |
| 10061-01-5 | cis-1,3-Dichloropropylene | 1.93 | 0.88 | U |
| 110-82-7 | Cyclohexane | 1.93 | 2.7 | |

Laboratory: York Analytical Laboratories, Inc. SDG: 22F1004
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Soil Vapor Laboratory ID: 22F1004-01 File ID: TQ220742.D
 Sampled: 06/16/22 13:00 Prepared: 06/23/22 02:00 Analyzed: 06/23/22 14:19
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BF21498 Sequence: Y2F2413 Calibration: YD20018 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 124-48-1 | Dibromochloromethane | 1.93 | 1.6 | U |
| 75-71-8 | Dichlorodifluoromethane | 1.93 | 3.1 | |
| 141-78-6 | Ethyl acetate | 1.93 | 1.4 | U |
| 100-41-4 | Ethyl Benzene | 1.93 | 17 | |
| 87-68-3 | Hexachlorobutadiene | 1.93 | 2.1 | UJ |
| 67-63-0 | Isopropanol | 1.93 | 1.4 | J- |
| 80-62-6 | Methyl Methacrylate | 1.93 | 0.79 | U |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 1.93 | 2.0 | |
| 75-09-2 | Methylene chloride | 1.93 | 8.7 | |
| 142-82-5 | n-Heptane | 1.93 | 7.1 | |
| 110-54-3 | n-Hexane | 1.93 | 4.9 | |
| 95-47-6 | o-Xylene | 1.93 | 60 | |
| 179601-23-1 | p- & m- Xylenes | 1.93 | 45 | |
| 622-96-8 | p-Ethyltoluene | 1.93 | 39 | |
| 115-07-1 | Propylene | 1.93 | 8.8 | |
| 100-42-5 | Styrene | 1.93 | 0.82 | U |
| 127-18-4 | Tetrachloroethylene | 1.93 | 12 | |
| 109-99-9 | Tetrahydrofuran | 1.93 | 1.1 | U |
| 108-88-3 | Toluene | 1.93 | 11 | |
| 156-60-5 | trans-1,2-Dichloroethylene | 1.93 | 0.77 | U |
| 10061-02-6 | trans-1,3-Dichloropropylene | 1.93 | 0.88 | U |
| 79-01-6 | Trichloroethylene | 1.93 | 0.26 | U |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 1.93 | 1.7 | |
| 108-05-4 | Vinyl acetate | 1.93 | 0.68 | U |
| 593-60-2 | Vinyl bromide | 1.93 | 0.85 | U |
| 75-01-4 | Vinyl Chloride | 1.93 | 0.25 | U |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| ISTD: 1,4-Difluorobenzene | 2361202 | 13.561 | 3144562 | 13.564 | |
| ISTD: d5-Chlorobenzene | 2275650 | 18.821 | 2743461 | 18.824 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. SDG: 22F1004
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Outdoor Ambient Air Laboratory ID: 22F1004-02 File ID: TQ220721.D
 Sampled: 06/17/22 15:00 Prepared: 06/22/22 03:00 Analyzed: 06/22/22 11:04
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BF21412 Sequence: Y2F2411 Calibration: YD20018 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|----|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.17 | 0.80 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.17 | 0.64 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.17 | 0.80 | U |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 1.17 | 0.89 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.17 | 0.64 | U |
| 75-34-3 | 1,1-Dichloroethane | 1.17 | 0.47 | U |
| 75-35-4 | 1,1-Dichloroethylene | 1.17 | 0.12 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.17 | 0.87 | UJ |
| 95-63-6 | 1,2,4-Trimethylbenzene | 1.17 | 0.57 | U |
| 106-93-4 | 1,2-Dibromoethane | 1.17 | 0.90 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.17 | 0.70 | U |
| 107-06-2 | 1,2-Dichloroethane | 1.17 | 0.47 | U |
| 78-87-5 | 1,2-Dichloropropane | 1.17 | 0.54 | U |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 1.17 | 0.82 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 1.17 | 0.57 | U |
| 106-99-0 | 1,3-Butadiene | 1.17 | 0.77 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.17 | 0.70 | U |
| 142-28-9 | 1,3-Dichloropropane | 1.17 | 0.54 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.17 | 0.70 | U |
| 123-91-1 | 1,4-Dioxane | 1.17 | 0.84 | U |
| 78-93-3 | 2-Butanone | 1.17 | 0.69 | |
| 591-78-6 | 2-Hexanone | 1.17 | 0.96 | UJ |
| 107-05-1 | 3-Chloropropene | 1.17 | 1.8 | U |
| 108-10-1 | 4-Methyl-2-pentanone | 1.17 | 0.48 | UJ |
| 67-64-1 | Acetone | 1.17 | 7.2 | |
| 107-13-1 | Acrylonitrile | 1.17 | 0.25 | U |
| 71-43-2 | Benzene | 1.17 | 0.56 | |
| 100-44-7 | Benzyl chloride | 1.17 | 0.60 | U |
| 75-27-4 | Bromodichloromethane | 1.17 | 0.78 | U |
| 75-25-2 | Bromoform | 1.17 | 1.2 | U |
| 74-83-9 | Bromomethane | 1.17 | 0.45 | U |
| 75-15-0 | Carbon disulfide | 1.17 | 0.36 | U |
| 56-23-5 | Carbon tetrachloride | 1.17 | 0.37 | |
| 108-90-7 | Chlorobenzene | 1.17 | 0.54 | U |
| 75-00-3 | Chloroethane | 1.17 | 0.31 | U |
| 67-66-3 | Chloroform | 1.17 | 0.57 | U |
| 74-87-3 | Chloromethane | 1.17 | 1.3 | |
| 156-59-2 | cis-1,2-Dichloroethylene | 1.17 | 0.12 | U |
| 10061-01-5 | cis-1,3-Dichloropropylene | 1.17 | 0.53 | U |
| 110-82-7 | Cyclohexane | 1.17 | 0.40 | U |

Laboratory: York Analytical Laboratories, Inc. SDG: 22F1004
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Outdoor Ambient Air Laboratory ID: 22F1004-02 File ID: TQ220721.D
 Sampled: 06/17/22 15:00 Prepared: 06/22/22 03:00 Analyzed: 06/22/22 11:04
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BF21412 Sequence: Y2F2411 Calibration: YD20018 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|----|
| 124-48-1 | Dibromochloromethane | 1.17 | 0.99 | U |
| 75-71-8 | Dichlorodifluoromethane | 1.17 | 2.8 | |
| 141-78-6 | Ethyl acetate | 1.17 | 0.97 | |
| 100-41-4 | Ethyl Benzene | 1.17 | 0.51 | U |
| 87-68-3 | Hexachlorobutadiene | 1.17 | 1.2 | UJ |
| 67-63-0 | Isopropanol | 1.17 | 1.8 | J- |
| 80-62-6 | Methyl Methacrylate | 1.17 | 0.48 | U |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 1.17 | 0.42 | U |
| 75-09-2 | Methylene chloride | 1.17 | 0.81 | U |
| 142-82-5 | n-Heptane | 1.17 | 0.48 | U |
| 110-54-3 | n-Hexane | 1.17 | 0.45 | |
| 95-47-6 | o-Xylene | 1.17 | 0.51 | |
| 179601-23-1 | p- & m- Xylenes | 1.17 | 1.4 | |
| 622-96-8 | p-Ethyltoluene | 1.17 | 0.57 | U |
| 115-07-1 | Propylene | 1.17 | 0.20 | U |
| 100-42-5 | Styrene | 1.17 | 0.50 | U |
| 127-18-4 | Tetrachloroethylene | 1.17 | 0.79 | U |
| 109-99-9 | Tetrahydrofuran | 1.17 | 0.69 | U |
| 108-88-3 | Toluene | 1.17 | 1.7 | |
| 156-60-5 | trans-1,2-Dichloroethylene | 1.17 | 0.46 | U |
| 10061-02-6 | trans-1,3-Dichloropropylene | 1.17 | 0.53 | U |
| 79-01-6 | Trichloroethylene | 1.17 | 0.16 | U |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 1.17 | 1.6 | |
| 108-05-4 | Vinyl acetate | 1.17 | 0.41 | U |
| 593-60-2 | Vinyl bromide | 1.17 | 0.51 | U |
| 75-01-4 | Vinyl Chloride | 1.17 | 0.15 | U |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 505547 | 11.972 | 679348 | 11.979 | |
| ISTD: 1,4-Difluorobenzene | 2839122 | 13.551 | 3753147 | 13.558 | |
| ISTD: d5-Chlorobenzene | 2436942 | 18.824 | 3277670 | 18.824 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. SDG: 22F1004
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Indoor Ambient Air Laboratory ID: 22F1004-03 File ID: TQ220723.D
 Sampled: 06/17/22 15:00 Prepared: 06/22/22 03:00 Analyzed: 06/22/22 13:23
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BF21412 Sequence: Y2F2411 Calibration: YD20018 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|------------|---|----------|----------------------------|---|
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 1.1 | 0.75 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 1.1 | 0.60 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 1.1 | 0.75 | U |
| 76-13-1 | 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 1.1 | 0.84 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 1.1 | 0.60 | U |
| 75-34-3 | 1,1-Dichloroethane | 1.1 | 0.44 | U |
| 75-35-4 | 1,1-Dichloroethylene | 1.1 | 0.11 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.1 | 0.81 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 1.1 | 7.5 | |
| 106-93-4 | 1,2-Dibromoethane | 1.1 | 0.84 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 1.1 | 0.66 | U |
| 107-06-2 | 1,2-Dichloroethane | 1.1 | 0.44 | U |
| 78-87-5 | 1,2-Dichloropropane | 1.1 | 0.51 | U |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane | 1.1 | 0.77 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 1.1 | 2.3 | |
| 106-99-0 | 1,3-Butadiene | 1.1 | 0.73 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 1.1 | 0.66 | U |
| 142-28-9 | 1,3-Dichloropropane | 1.1 | 0.51 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 1.1 | 1.2 | |
| 123-91-1 | 1,4-Dioxane | 1.1 | 0.79 | U |
| 78-93-3 | 2-Butanone | 1.1 | 2.3 | |
| 591-78-6 | 2-Hexanone | 1.1 | 0.90 | U |
| 107-05-1 | 3-Chloropropene | 1.1 | 1.7 | U |
| 108-10-1 | 4-Methyl-2-pentanone | 1.1 | 0.45 | U |
| 107-13-1 | Acrylonitrile | 1.1 | 0.24 | U |
| 71-43-2 | Benzene | 1.1 | 0.63 | |
| 100-44-7 | Benzyl chloride | 1.1 | 0.57 | U |
| 75-27-4 | Bromodichloromethane | 1.1 | 0.73 | U |
| 75-25-2 | Bromoform | 1.1 | 1.1 | U |
| 74-83-9 | Bromomethane | 1.1 | 0.43 | U |
| 75-15-0 | Carbon disulfide | 1.1 | 0.34 | U |
| 56-23-5 | Carbon tetrachloride | 1.1 | 0.28 | |
| 108-90-7 | Chlorobenzene | 1.1 | 0.50 | U |
| 75-00-3 | Chloroethane | 1.1 | 0.29 | U |
| 67-66-3 | Chloroform | 1.1 | 0.53 | U |
| 74-87-3 | Chloromethane | 1.1 | 1.2 | |
| 156-59-2 | cis-1,2-Dichloroethylene | 1.1 | 0.11 | U |
| 10061-01-5 | cis-1,3-Dichloropropylene | 1.1 | 0.50 | U |
| 110-82-7 | Cyclohexane | 1.1 | 0.41 | |
| 124-48-1 | Dibromochloromethane | 1.1 | 0.93 | U |

Laboratory: York Analytical Laboratories, Inc. SDG: 22F1004
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Indoor Ambient Air Laboratory ID: 22F1004-03 File ID: TQ220723.D
 Sampled: 06/17/22 15:00 Prepared: 06/22/22 03:00 Analyzed: 06/22/22 13:23
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BF21412 Sequence: Y2F2411 Calibration: YD20018 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|-------------|-----------------------------------|----------|----------------------------|---|
| 75-71-8 | Dichlorodifluoromethane | 1.1 | 2.5 | |
| 141-78-6 | Ethyl acetate | 1.1 | 2.7 | |
| 100-41-4 | Ethyl Benzene | 1.1 | 83 | |
| 87-68-3 | Hexachlorobutadiene | 1.1 | 1.2 | U |
| 67-63-0 | Isopropanol | 1.1 | 23 | |
| 80-62-6 | Methyl Methacrylate | 1.1 | 6.2 | |
| 1634-04-4 | Methyl tert-butyl ether (MTBE) | 1.1 | 0.39 | U |
| 75-09-2 | Methylene chloride | 1.1 | 0.76 | U |
| 142-82-5 | n-Heptane | 1.1 | 1.3 | |
| 110-54-3 | n-Hexane | 1.1 | 1.1 | |
| 95-47-6 | o-Xylene | 1.1 | 75 | |
| 179601-23-1 | p- & m- Xylenes | 1.1 | 280 | |
| 622-96-8 | p-Ethyltoluene | 1.1 | 7.5 | |
| 115-07-1 | Propylene | 1.1 | 0.19 | U |
| 100-42-5 | Styrene | 1.1 | 0.47 | U |
| 127-18-4 | Tetrachloroethylene | 1.1 | 0.74 | U |
| 109-99-9 | Tetrahydrofuran | 1.1 | 0.65 | U |
| 108-88-3 | Toluene | 1.1 | 2.3 | |
| 156-60-5 | trans-1,2-Dichloroethylene | 1.1 | 0.43 | U |
| 10061-02-6 | trans-1,3-Dichloropropylene | 1.1 | 0.50 | U |
| 79-01-6 | Trichloroethylene | 1.1 | 0.15 | U |
| 75-69-4 | Trichlorofluoromethane (Freon 11) | 1.1 | 1.4 | |
| 108-05-4 | Vinyl acetate | 1.1 | 0.39 | U |
| 593-60-2 | Vinyl bromide | 1.1 | 0.48 | U |
| 75-01-4 | Vinyl Chloride | 1.1 | 0.14 | U |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 496456 | 12.005 | 679348 | 11.979 | |
| ISTD: 1,4-Difluorobenzene | 2652947 | 13.574 | 3753147 | 13.558 | |
| ISTD: d5-Chlorobenzene | 2473014 | 18.824 | 3277670 | 18.824 | |

* Values outside of QC limits

JR 12/8/2025

Laboratory: York Analytical Laboratories, Inc. SDG: 22F1004
 Client: Advanced Cleanup Technologies, Inc. Project: 9628-BXNY
 Matrix: Indoor Ambient Air Laboratory ID: 22F1004-03RE1 File ID: TQ220737.D
 Sampled: 06/17/22 15:00 Prepared: 06/23/22 02:00 Analyzed: 06/23/22 09:05
 Solids: Preparation: EPA TO15 PREP Initial/Final: 400 mL / 400 mL
 Batch: BF21498 Sequence: Y2F2413 Calibration: YD20018 Instrument: TO15 AIR2

| CAS NO. | COMPOUND | DILUTION | CONC. (ug/m ³) | Q |
|---------|----------|----------|----------------------------|---|
| 67-64-1 | Acetone | 41.1 | 3300 | |

| INTERNAL STANDARD | AREA | RT | REF AREA | REF RT | Q |
|---------------------------|---------|--------|----------|--------|---|
| Bromochloromethane | 391726 | 11.979 | 554760 | 11.985 | |
| ISTD: 1,4-Difluorobenzene | 2170457 | 13.558 | 3144562 | 13.564 | |
| ISTD: d5-Chlorobenzene | 1751490 | 18.821 | 2743461 | 18.824 | |

* Values outside of QC limits

JR 12/8/2025

December 8, 2025

Mr. Jason Stewart
Advanced Cleanup Technologies
228 Park Ave S PMB 34864
New York, New York 10003

Re: Data Usability Summary Report – Phoenix Environmental Laboratories, Inc. - GCL97021 V1

Dear Mr. Stewart:

The evaluation of analytical data by Phoenix Environmental Laboratories for project 9268-BXNY, which were reported in a single data package under Job No. GCL97021 V1 has been completed. The following samples were reported.

IA-4 IA-7

Analysis was performed in accordance with EPA Method TO-15 (volatile organics). The review was performed to the extent possible, in accordance with the analytical method, and “DER-10/ Technical Guidance for Site Investigation and Remediation”. Professional judgment is applied as necessary and appropriate. National Functional Guidelines for Organic Data Review was consulted as needed. Qualifiers consistent with those defined by EPA Region 2 are applied as necessary and appropriate.

Below is the Data Usability Summary Report (DUSR) associated with these samples.

| Data Usability Summary Report | |
|---|--------------------------------|
| 1. Is the data package complete as defined under the requirements for the most current DEC ASP Category B or USEPA CLP data deliverables? | No – see Documentation Section |
| 2. Have all holding times been met? | Yes |
| 3. Do all the QC data; blanks, instrument tunings, calibration standards, calibration verifications, surrogate recoveries, spike recoveries, replicate analyses, laboratory controls and sample data fall within the protocol required limits and specifications? | No -see following sections |
| 4. Have all of the data been generated using established and agreed upon analytical protocols? | Yes |
| 5. Does an evaluation of the raw data confirm the results provided in the data summary sheet and the quality control verification forms? | Yes |
| 6. Have the correct data qualifiers been used and are they consistent with the most current DEC ASP? | Yes |

| | |
|---|-----|
| 7. Have any quality control (QC) exceedances been specifically noted in the DUSR and have the corresponding QC summary sheet from the data package been attached to the DUSR? | Yes |
|---|-----|

Overall Evaluation

Based on the data review effort, results are usable, with the following qualifications. For samples that are qualified as estimated (J-, UJ), detected results may be biased low. False negatives may exist in non-detect results. Sample results that are qualified as estimated (J+) may be biased high. For samples that are qualified as estimated with any combination of (J), (J-) and/or (J+), the (J) qualifier takes precedence and is applied to the sample result. It is not possible to determine the direction of the bias and the overall effect on the result. Where a sample result is rejected (R), the 'R' qualifier takes precedence over any other qualification.

- The non-detect results for benzyl chloride, n-butylbenzene, 1,2,4-trichlorobenzene and hexachlorobutadiene in IA-4 and IA-7 are rejected (R) due to unacceptable initial calibration.
- The results for ethanol, 1,3-dichlorobenzene, 1,2-dichlorobenzene, 1,2,4-trichlorobenzene and hexachlorobutadiene in IA-4 and IA-7 are qualified as estimated (J) due to unacceptable recovery in the second source initial calibration verification (ICV) standard.
- The results for chloromethane, ethanol and acetone in IA-1 and IA-7 are qualified as estimated (J+) due to high recovery in the LCS.
- The results for cyclohexane in IA-4 and IA-7 are qualified as not detected (U) at the reported value due to poor spectral match.
- The results for ethanol in IA-4 and acetone in IA-7 are qualified as estimated (J) because the reported concentration is outside the calibration range.

Qualifier definitions are provided in Attachment A. A copy of the chain of custody record is provided in Attachment B. Pages from the data package illustrating the exceedances and issues described in this validation report are included in Attachment C. Annotated Summary Forms are included in Attachment D detailing qualifications resulting from the data review effort.

The following components were reviewed, where applicable:

- Chain of Custody
- Receiving conditions
- Holding times
- Preservation

- Analyte lists
- Reporting limits
- Requested methods
- Units, and
- Sample related quality control data:
 - Method, instrument blanks
 - Clean canister certification
 - Field blanks
 - Surrogate recoveries
 - LCS recoveries
 - Internal standard area response
 - Duplicates
 - Analyte Identification
- Instrument related quality control data:
 - Instrument tunes
 - Calibration summaries

The following sections of the report detail only quality control exceedances that impacted results. Where a quality control item exceeded control limits but there is no impact to the samples results, these are not detailed in the report.

Documentation: A completeness review of the data package was performed, and the data package was determined to be a complete Category B data package with the following exceptions:

- The data package does not include summary forms and raw data for the second source initial calibration verification (ICV) standard. The laboratory was contacted and provided the ICV data and is included in attachment C.
- Calibration data for the clean canister certification analysis is not included in the data package.

The following documentation issues were observed during the review:

- Improper edits were observed on the chain of custody records. When edits are made, a single line should be drawn through the error, then dated and initialed by the person making the edits.
- Results for ethanol, acetone, isopropyl alcohol, and hexane in IA-4 and IA-7 are flagged 'S' which the data deliverable identifies this flag as 'This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level.' However, all sample detections are greater than five times the reporting limit. The validator removed the 'S' qualifier.
- The data package documents a sample collection date of August 2, 2022; however, the samples were collected on August 1, 2022.

Holding Times, Preservation, Sample Integrity:

A copy of the applicable chain of custody (COC) record was included in the data package, documenting sample collection date of August 1, 2022. The samples were received on August 3, 2022. Samples were received intact and were analyzed within the method specified hold time.

A. Volatile Organics**1. Calibration**

One initial calibration (IC) was performed in support of the sample analysis. All relative response factors (RRFs) and relative standard deviations (RSDs) or correlation coefficients are acceptable with the exceptions noted below.

| Analyte | RSD |
|------------------------------|-----|
| Benzyl chloride (SIM) | 167 |
| n-Butylbenzene (SIM) | 129 |
| 1,2,4-Trichlorobenzene (SIM) | 173 |
| Hexachlorobutadiene (SIM) | 167 |

The laboratory applied quadratic equations to achieve an acceptable correlation coefficient of 1.00. However, a review of the individual calibration levels indicates a lack of instrument sensitivity in the lower calibration standards as presented below.

| Analyte | RRF 0.01 | RRF 0.02 | RRF 0.035 | RRF 0.05 | RRF 0.1 | RRF 0.2 | RRF 0.5 | RRF 1.0 | RRF 2.5 | RRF 5.0 |
|------------------------------|------------------|------------------|--------------|--------------|------------|------------|------------|------------|------------|------------|
| Benzyl chloride (SIM) | 0.01 2 | 0.01 7 | 0.028 | 0.050 | 0.111 | 0.169 | 0.462 | 1.088 | 2.477 | - |
| n-Butylbenzene (SIM) | - | - | - | 0.034 | 0.114 | 0.192 | 0.479 | 1.044 | 2.487 | - |
| 1,2,4-Trichlorobenzene (SIM) | 0.03 6 | 0.03 7 | 0.038 | 0.040 | 0.098 | 0.164 | 0.393 | 1.216 | 2.348 | 5.045 |
| Hexachlorobutadiene (SIM) | 0.01 6 | 0.02 2 | 0.032 | 0.042 | 0.116 | 0.200 | 0.483 | 1.002 | 2.502 | - |

Removing the standards with low responses (bolded above) still produces an RSD >30% as follows:

Benzyl chloride (SIM) -114 RSD
 n-Butylbenzene (SIM) -113 RSD
 1,2,4-Trichlorobenzene (SIM) – 124 RSD
 Hexachlorobutadiene (SIM) – 114RSD

Based on professional judgment the non-detect results for benzyl chloride, n-butylbenzene, 1,2,4-trichlorobenzene and hexachlorobutadiene in IA-4 and IA-7 are rejected (R) due to unacceptable initial calibration.

A second source ICV standard was analyzed after the IC, and all percent differences are acceptable (70-130%R), with the following exceptions.

| Analyte | %R | Associated Sample | Qualifier Applied |
|-----------------------------------|------|-------------------|-------------------|
| <i>ICV CHEM20 7/20/2022 12:30</i> | | | |
| Ethanol | 174 | IA-4 | J, UJ |
| 1,3-Dichlorobenzene (SIM) | 136* | IA7 | |
| 1,2-Dichlorobenzene (SIM) | 139* | | |
| 1,2,4-Trichlorobenzene (SIM) | 160* | | |
| Hexachlorobutadiene (SIM) | NR* | | |

NR=not reported

*The summary form provided by the laboratory reports ICV recoveries based on the full scan analysis. The validator reviewed the raw data to determine the recovery in the SIM analysis and this recovery is presented above.

The results for ethanol, 1,3-dichlorobenzene, 1,2-dichlorobenzene, 1,2,4-trichlorobenzene and hexachlorobutadiene in IA-4 and IA-7 are qualified as estimated (UJ) due to unacceptable recovery in the ICV.

2. Laboratory Control Sample (LCS)

One LCS was prepared and analyzed with the sample. The LCS is evaluated using control limit of 70-130%R. All recoveries are acceptable, with the following exceptions:

| Analyte | LCS %R | Affected Sample | Qualifier Applied |
|---------------|--------|-----------------|-------------------|
| Chloromethane | 134 | IA-4 | J+ |
| Ethanol | 149 | IA-7 | |
| Acetone | 132 | | |

The results for chloromethane, ethanol and acetone in IA-1 and IA-7 are qualified as estimated (J+) due to high recovery in the LCS.

3. Compound Identification

The results for cyclohexane in IA-4 and IA-7 are qualified as not detected (U) at the reported value due to poor spectral match.

4. Compound Quantitation

The results for ethanol in IA-4 and acetone in IA-7 are qualified as estimated (J) because the reported concentration is outside the calibration range. Results from the diluted analysis should be used.

5. Clean Canister Certification

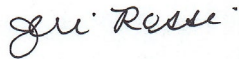
Ethanol is detected in one of the clean canister certification batches as presented below.

| Batch Canister | Analyte | Conc. (ppbv) |
|----------------|---------|--------------|
| Batch #259 | Ethanol | 0.739 |

The data package does not indicate which canisters are associated with which batch canister certification data. Since the concentration of ethanol in all samples is more than ten times the concentration in the clean canister certification no action is warranted.

No other sample results are qualified. Please feel free to contact me at (908) 370-3431 or richjerirossi513@gmail.com if you have any questions regarding this data package review report or need further information.

Sincerely,



Jeri L Rossi, CEAC

Environmental Consulting Chemist

ATTACHMENT A

Qualifier Definitions

EPA Qualifier Definitions

- U The analyte was analyzed for but was not detected above the level of the reported sample quantitation limit.
- J The result is an estimated quantity. The associated numerical value is the approximate concentration of the analyte in the sample.
- J+ The result is an estimated quantity, but the result may be biased high.
- J- The result is an estimated quantity, but the result may be biased low.
- NJ The analyte has been “tentatively identified” or “presumptively” as present and the associated numerical value is the estimated concentration in the sample.
- UJ The analyte was analyzed for but was not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
- R The data are unusable. The sample results are rejected due to serious deficiencies in meeting QC criteria. The analyte may or may not be present in the sample.

ATTACHMENT B

CHAIN OF CUSTODY (COC)



587 East Middle Turnpike, P.O. Box 370, Manchester, CT 06040
Telephone: (860) 645-1102 • Fax: (860) 645-3623

CHAIN OF CUSTODY RECORD

AIR ANALYSES

860-645-1102

email: greg@phoenixlabs.com

P.O. #

Page 1 of 1

Data Delivery:

Fax #: _____
 Email: _____
 Phone #: _____

ACT

| | | |
|--|--|--|
| Report to: jason Stewart | Project Name: q268 - Bx NY | Data Format: (Circle) Equis Excel Other: _____ |
| Customer: Advanced Cleanup Technologies, Inc. | Invoice to: Karen Friedman | Requested Deliverable: RCP ASP CAT B |
| Address: 228 Park Ave S PMB 34864 | | MCP NJ Deliverables |
| 13763 New York, New York 10003 | Sampled by: Yisng Yeng / Ray Buzeta | Quote Number: _____ |

| Phoenix ID # | Client Sample ID | Canister ID # | Canister Size (L) | Outgoing Canister Pressure ("Hg) | Incoming Canister Pressure ("Hg) | Flow Regulator ID # | Flow Controller Setting (mL/min) | Sampling Start Time | Sampling End Time | Sample Start Date | Canister Pressure at Start ("Hg) | Canister Pressure at End ("Hg) | Ambient/Indoor Air | Soil Gas | Grab (G) Composite (C) | TO-15 | APH |
|-------------------------------|------------------|---------------|-------------------|----------------------------------|----------------------------------|---------------------|----------------------------------|---------------------|-------------------|-------------------|----------------------------------|--------------------------------|--------------------|----------|------------------------|-------|-----|
| THIS SECTION FOR LAB USE ONLY | | | | | | | | | | | | | | | | | |
| 92021 | IA-4 | 28560 | 6.0L | -30 | 40 | 5043 | 3.6 | 15:00 8/1/22 | 11:00 8/2/22 | 8/1/22 | -30 | -8 | ✓ | | | ✓ | |
| 92022 | IA-7 | 23330 | 6.0L | -30 | 40 5 | 5000 | 3.8 | 15:00 8/1/22 | 10:22 8/2/22 | 8/1/22 | -30 | -6 | ✓ | | | ✓ | |

| | | | | |
|---|---------------------------------|--------------------------------------|-----------------------------------|---|
| Relinquished by: ACT <i>Ray Buzeta</i> 8/2/22 3:15 PM | Accepted by: <i>[Signature]</i> | Date: 8-3-22 8/3/22 | Time: 10:54 1750 | I attest that all media released by Phoenix Environmental Laboratories, Inc. have been received in good working condition and agree to the terms and conditions as listed on the back of this document. |
| Signature: _____ | | | Date: _____ | |

| | | | | | | | |
|---|--|--|---|---|------------------------------|---|--|
| State Where Samples Collected: _____ | Turnaround Time: 1 Day <input type="checkbox"/> 2 Day <input type="checkbox"/> 3 Day <input type="checkbox"/> 4 Day <input type="checkbox"/> 5 Day <input type="checkbox"/> | Requested Criteria: (Please Circle) CT: _____ TAC I/C TAC RES SVVC I/C SVVC RES GWV I/C GWV CES | MA: _____ Indoor Air: Residential Ind/Commercial Soil Gas: Residential Ind/Commercial | NI: _____ Indoor Air: Residential Ind/Commercial Soil Gas: Residential Ind/Commercial | NY: _____ Vapor Intrusion | PA: _____ Indoor Air: Residential Non-residential | VT: _____ Indoor Air: Residential Industrial Sub-slab Residential Industrial |
| SPECIAL INSTRUCTIONS, OC REQUIREMENTS, REGULATORY INFORMATION: Can # 23338 (3) - 6.0L 24 hr returned unused | | | | | | | |

01/09/2024

Phoenix Environmental Laboratories, Inc.

Page 19 of 195

ATTACHMENT C

**SELECTED PAGES FROM DATA PACKAGE –
QC EXCEEDANCES AND VALIDATION ISSUES**

3
AIR LCS RECOVERY

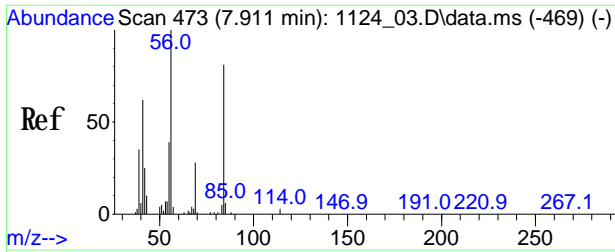
Lab Name: Phoenix Environmental Labs Client: ACT

Lab Code: Phoenix Case No: _____ SAS No: _____ SDG No GCL97021

LCS - Client Id: CL97022 LCS

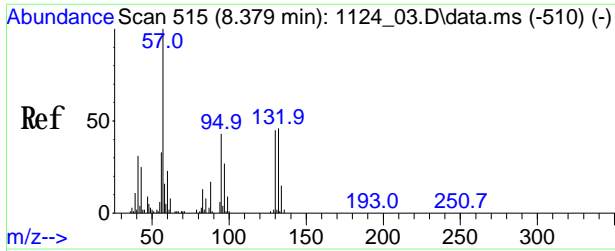
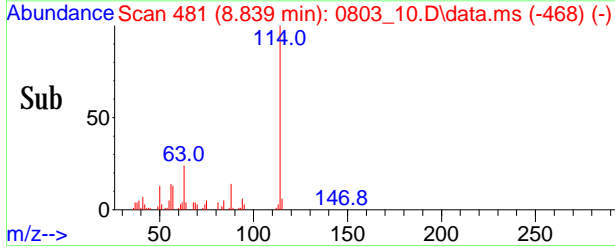
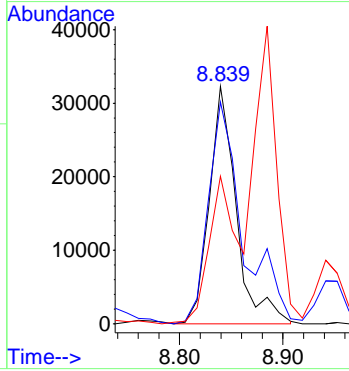
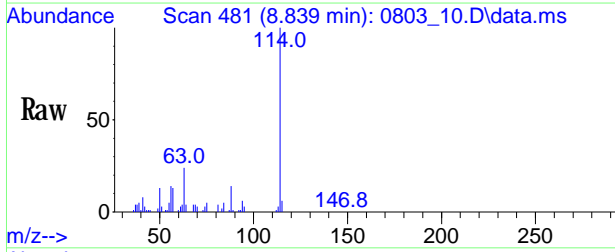
| COMPOUND | SPIKE ADDED (ppbv) | | LCS CONCENTRATION (ppbv) | LCS % REC # | QC. LIMITS REC. | |
|-------------------------------|--------------------|--|--------------------------|-------------|-----------------|-----|
| Propylene | 10 | | 9.838 | 98 | 70 | 130 |
| Dichlorodifluoromethane | 10 | | 10.74 | 107 | 70 | 130 |
| Chloromethane | 10 | | 13.42 | 134 * | 70 | 130 |
| 1,2-Dichlorotetrafluoroethane | 10 | | 10.73 | 107 | 70 | 130 |
| Vinyl Chloride | 10 | | 10.76 | 108 | 70 | 130 |
| 1,3-Butadiene | 10 | | 11.96 | 120 | 70 | 130 |
| Bromomethane | 10 | | 9.751 | 98 | 70 | 130 |
| Chloroethane | 10 | | 9.310 | 93 | 70 | 130 |
| Ethanol | 10 | | 14.93 | 149 * | 70 | 130 |
| Acetone | 10 | | 13.17 | 132 * | 70 | 130 |
| Trichlorofluoromethane | 10 | | 10.69 | 107 | 70 | 130 |
| Isopropylalcohol | 10 | | 13.04 | 130 | 70 | 130 |
| Acrylonitrile | 10 | | 12.93 | 129 | 70 | 130 |
| 1,1-Dichloroethene | 10 | | 11.03 | 110 | 70 | 130 |
| Methylene Chloride | 10 | | 12.51 | 125 | 70 | 130 |
| Carbon Disulfide | 10 | | 10.54 | 105 | 70 | 130 |
| Trichlorotrifluoroethane | 10 | | 10.75 | 108 | 70 | 130 |
| Trans-1,2-Dichloroethene | 10 | | 10.80 | 108 | 70 | 130 |
| 1,1-Dichloroethane | 10 | | 10.62 | 106 | 70 | 130 |
| Methyl tert-butyl ether(MTBE) | 10 | | 9.709 | 97 | 70 | 130 |
| Methyl Ethyl Ketone | 10 | | 12.90 | 129 | 70 | 130 |
| Cis-1,2-Dichloroethene | 10 | | 11.69 | 117 | 70 | 130 |
| Hexane | 10 | | 11.06 | 111 | 70 | 130 |
| Chloroform | 10 | | 10.20 | 102 | 70 | 130 |
| Ethyl acetate | 10 | | 9.384 | 94 | 70 | 130 |
| Tetrahydrofuran | 10 | | 12.76 | 128 | 70 | 130 |
| 1,2-Dichloroethane | 10 | | 10.95 | 110 | 70 | 130 |
| 1,1,1-Trichloroethane | 10 | | 10.36 | 104 | 70 | 130 |
| Benzene | 10 | | 9.791 | 98 | 70 | 130 |
| Carbon Tetrachloride | 10 | | 11.21 | 112 | 70 | 130 |
| Cyclohexane | 10 | | 9.883 | 99 | 70 | 130 |
| 1,2-dichloropropane | 10 | | 10.35 | 104 | 70 | 130 |
| Bromodichloromethane | 10 | | 11.14 | 111 | 70 | 130 |
| Trichloroethene | 10 | | 10.20 | 102 | 70 | 130 |
| 1,4-Dioxane | 10 | | 9.620 | 96 | 70 | 130 |
| Heptane | 10 | | 12.61 | 126 | 70 | 130 |
| cis-1,3-Dichloropropene | 10 | | 10.24 | 102 | 70 | 130 |
| 4-Methyl-2-pentanone(MIBK) | 10 | | 12.75 | 128 | 70 | 130 |
| trans-1,3-Dichloropropene | 10 | | 10.22 | 102 | 70 | 130 |
| 1,1,2-Trichloroethane | 10 | | 10.04 | 100 | 70 | 130 |
| Toluene | 10 | | 10.00 | 100 | 70 | 130 |
| Dibromochloromethane | 10 | | 10.98 | 110 | 70 | 130 |
| 2-Hexanone(MBK) | 10 | | 12.78 | 128 | 70 | 130 |
| 1,2-Dibromoethane(EDB) | 10 | | 10.16 | 102 | 70 | 130 |

FORM III AIR



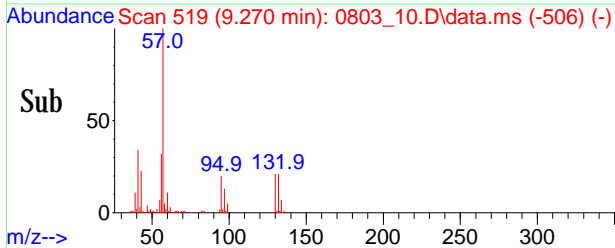
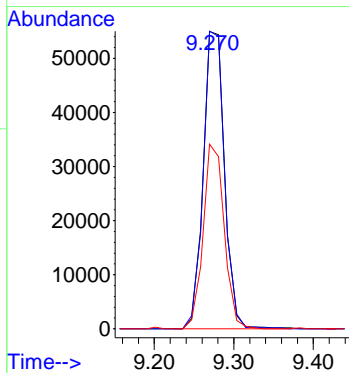
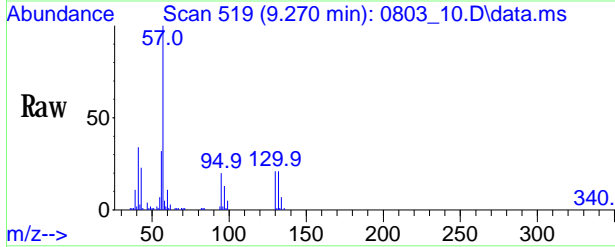
#36
 Cyclohexane
 Conc: 8S 1.804 ppbv
 RT: 8.839 min Scan# 481
 Delta R.T. 0.000 min
 Lab File: 0803_10.D
 Acq: 3 Aug 2022 10:55 pm

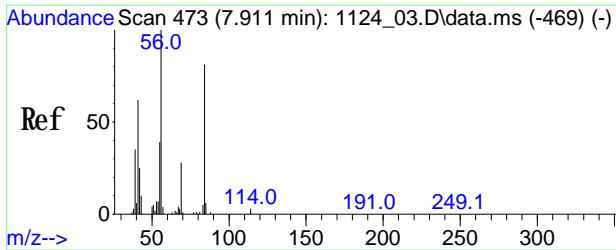
| Tgt Ion | Ratio | Resp | Lower | Upper |
|---------|-------|-------|-------|-------|
| 84 | 100 | 58496 | | |
| 55 | 119.9 | 41.2 | 61.8# | |
| 42 | 64.4 | 28.9 | 43.3# | |



#40
 Trichloroethene
 Conc: 8S 2.299 ppbv
 RT: 9.270 min Scan# 519
 Delta R.T. -0.000 min
 Lab File: 0803_10.D
 Acq: 3 Aug 2022 10:55 pm

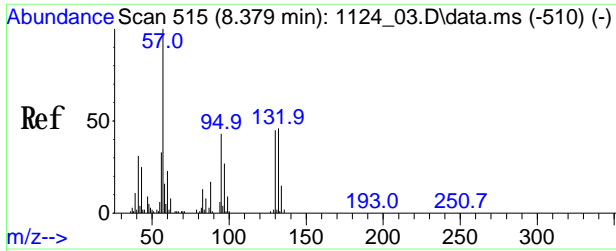
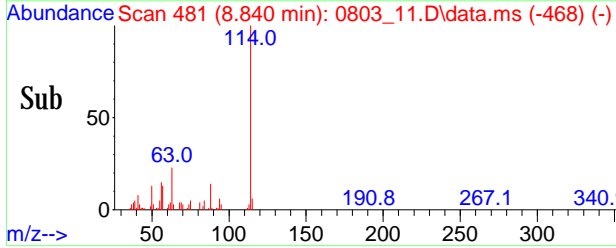
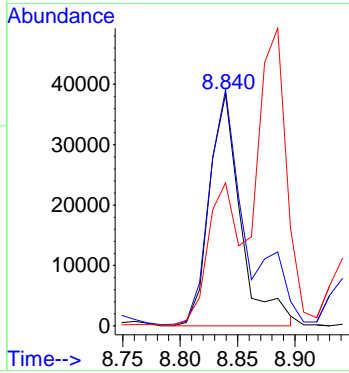
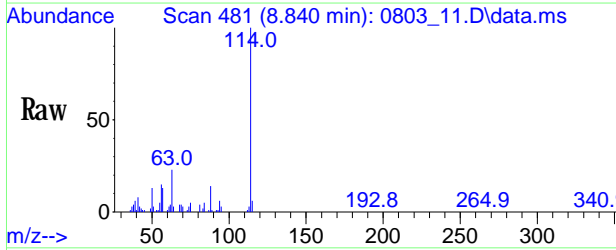
| Tgt Ion | Ratio | Resp | Lower | Upper |
|---------|-------|--------|-------|-------|
| 130 | 100 | 102505 | | |
| 132 | 99.4 | 79.9 | 119.9 | |
| 97 | 61.3 | 52.9 | 79.3 | |





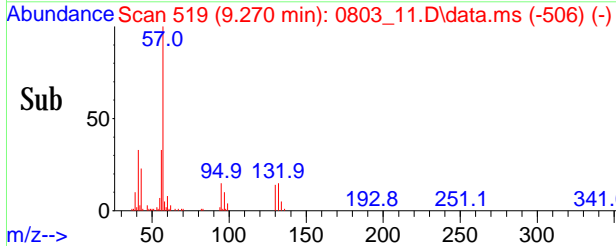
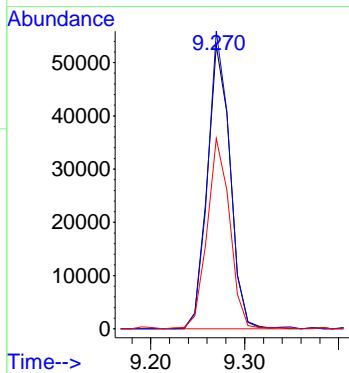
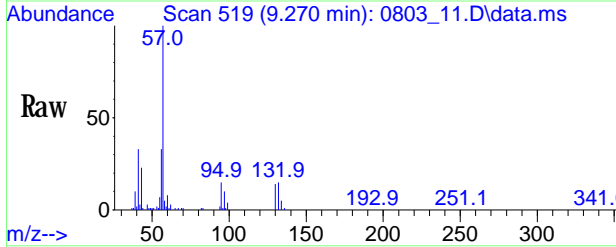
#36
 Cyclohexane
 Conc: 8S 2.159 ppbv
 RT: 8.840 min Scan# 481
 Delta R.T. 0.001 min
 Lab File: 0803_11.D
 Acq: 3 Aug 2022 11:31 pm

Tgt Ion: 84 Resp: 72982
 Ion Ratio Lower Upper
 84 100
 55 120.3 41.2 61.8#
 42 56.7 28.9 43.3#



#40
 Trichloroethene
 Conc: 8S 1.895 ppbv
 RT: 9.270 min Scan# 519
 Delta R.T. -0.000 min
 Lab File: 0803_11.D
 Acq: 3 Aug 2022 11:31 pm

Tgt Ion: 130 Resp: 89980
 Ion Ratio Lower Upper
 130 100
 132 100.8 79.9 119.9
 97 66.5 52.9 79.3



1
AIR ANALYSIS DATA SHEET

CLIENT ID

CANISTER BLK 2599

| | | | |
|--------------|----------|------------------|----------------------|
| Client: | ACT | Lab: | Phoenix Env. Labs |
| SDG No.: | GCL97021 | Lab Sample ID: | CANISTER BLK 2599 |
| Canister: | CANBL | Lab File ID: | 0713_08.D |
| Instrument: | CHEM20 | Column: | _____ |
| Purge Volume | 200 | (cc) | Date Received: _____ |
| Matrix: | AIR | Dilution Factor: | 1 |
| | | Date Analyzed: | 07/13/22 |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|------------|------------------------------------|-------|---|-------|-------|---|
| 115-07-1 | Propylene | 0.581 | U | 0.581 | 0.581 | r |
| 75-71-8 | Dichlorodifluoromethane | 0.202 | U | 0.202 | 0.202 | r |
| 74-87-3 | Chloromethane | 0.485 | U | 0.485 | 0.485 | r |
| 106-99-0 | 1,3-Butadiene | 0.452 | U | 0.452 | 0.452 | r |
| 75-00-3 | Chloroethane | 0.379 | U | 0.379 | 0.379 | r |
| 64-17-5 | Ethanol | 0.739 | | 0.531 | 0.531 | r |
| 67-64-1 | Acetone | 0.421 | U | 0.421 | 0.421 | r |
| 67-63-0 | Isopropylalcohol | 0.407 | U | 0.407 | 0.407 | r |
| 107-13-1 | Acrylonitrile | 0.461 | U | 0.461 | 0.461 | r |
| 75-09-2 | Methylene Chloride | 0.863 | U | 0.863 | 0.863 | r |
| 75-15-0 | Carbon Disulfide | 0.321 | U | 0.321 | 0.321 | r |
| 1634-04-4 | Methyl tert-butyl ether(MTBE) | 0.278 | U | 0.278 | 0.278 | r |
| 78-93-3 | Methyl Ethyl Ketone | 0.339 | U | 0.339 | 0.339 | r |
| 110-54-3 | Hexane | 0.284 | U | 0.284 | 0.284 | r |
| 141-78-6 | Ethyl acetate | 0.278 | U | 0.278 | 0.278 | r |
| 109-99-9 | Tetrahydrofuran | 0.339 | U | 0.339 | 0.339 | r |
| 110-82-7 | Cyclohexane | 0.291 | U | 0.291 | 0.291 | r |
| 142-82-5 | Heptane | 0.244 | U | 0.244 | 0.244 | r |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | 0.244 | U | 0.244 | 0.244 | r |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.220 | U | 0.220 | 0.220 | r |
| 108-88-3 | Toluene | 0.266 | U | 0.266 | 0.266 | r |
| 591-78-6 | 2-Hexanone(MBK) | 0.244 | U | 0.244 | 0.244 | r |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.146 | U | 0.146 | 0.146 | r |
| 108-90-7 | Chlorobenzene | 0.217 | U | 0.217 | 0.217 | r |
| 100-41-4 | Ethylbenzene | 0.230 | U | 0.230 | 0.230 | r |
| 100-42-5 | Styrene | 0.235 | U | 0.235 | 0.235 | r |
| 95-47-6 | o-Xylene | 0.230 | U | 0.230 | 0.230 | r |
| 98-82-8 | Isopropylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 622-96-8 | 4-Ethyltoluene | 0.204 | U | 0.204 | 0.204 | r |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane(sim) | 0.143 | U | 0.143 | 0.143 | r |
| 75-01-4 | Vinyl Chloride(sim) | 0.078 | U | 0.078 | 0.078 | r |
| 74-83-9 | Bromomethane(sim) | 0.258 | U | 0.258 | 0.258 | r |
| 75-69-4 | Trichlorofluoromethane(sim) | 0.178 | U | 0.178 | 0.178 | r |
| 107-06-2 | 1,2-Dichloroethane(sim) | 0.247 | U | 0.247 | 0.247 | r |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

6B
AIR INITIAL CALIBRATION DATA

Lab Name: Phoenix Environmental Labs
 Lab Code: Phoenix
 Instrument ID: CHEM20
 Heated Purge (Y/N): Y
 GC Column: _____

Client: _____
 SDG No.: GCL97021
 Calibration Date From: 07/19/22 20:00
 Calibration Date Thru: 07/19/22 23:55
 Method File: 20_AIR_0719.M

Laboratory File Ids

| | | | | | | | | | | | |
|-------|------------------|------|------------------|------|------------------|-------|------------------|-------|------------------|-------|------------------|
| RRF1 | <u>0719_05.D</u> | RRF2 | <u>0719_06.D</u> | RRF3 | <u>0719_07.D</u> | RRF4 | <u>0719_08.D</u> | RRF5 | <u>0719_09.D</u> | RRF6 | <u>0719_10.D</u> |
| RRF7 | <u>0719_11.D</u> | RRF8 | <u>0719_16.D</u> | RRF9 | <u>0719_12.D</u> | RRF10 | <u>0719_13.D</u> | RRF11 | <u>0719_17.D</u> | RRF12 | <u>0719_14.D</u> |
| RRF13 | <u>0719_15.D</u> | | | | | | | | | | |

| COMPOUND | RRF1 0.01 | RRF2 0.02 | RRF3 0.035 | RRF4 0.05 | RRF5 0.1 | RRF6 0.2 | RRF7 0.5 | RRF8 1 | RRF9 2.5 | RRF10 5 | RRF11 10 | RRF12 25 | RRF13 40 | RRF | % RSD |
|--------------------------------|--------------|--------------|---------------|--------------|-------------|-------------|-------------|-----------|-------------|------------|-------------|-------------|-------------|---------|----------|
| Benzene(sim) | 3.743 | 3.422 | 3.408 | 2.815 | 3.175 | 2.837 | 2.526 | 2.534 | | | | | | 3.057 | 14.65 |
| Carbon Tetrachloride(sim) | 2.056 | 1.918 | 1.872 | 1.719 | 2.431 | 2.142 | 2.007 | 2.130 | 2.031 | | | | | 2.034 | 9.82 |
| 1,1-Dichloroethene(sim) | 2.029 | 1.838 | 1.664 | 1.569 | 2.147 | 1.872 | 1.803 | 1.776 | 1.680 | 1.880 | | | | 1.826 | 9.42 |
| Trichlorotrifluoroethane(sim) | 2.032 | 1.865 | 1.755 | 1.832 | 2.421 | 2.141 | 2.012 | 2.024 | 1.968 | 2.164 | | | | 2.021 | 9.46 |
| Trans-1,2-Dichloroethene(sim) | 1.88 | 1.586 | 1.546 | 1.275 | 2.016 | 1.703 | 1.539 | 1.547 | 1.537 | 1.736 | | | | 1.637 | 12.66 |
| 1,1-Dichloroethane(sim) | 1.956 | 1.906 | 1.825 | 1.800 | 2.345 | 2.106 | 1.959 | 1.978 | 1.950 | 2.174 | | | | 2.000 | 8.28 |
| Cis-1,2-Dichloroethene(sim) | 1.724 | 1.626 | 1.519 | 1.165 | 1.768 | 1.662 | 1.600 | 1.528 | 1.477 | 1.675 | | | | 1.575 | 10.87 |
| Chloroform(sim) | 2.019 | 1.659 | 1.745 | 1.688 | 2.293 | 2.011 | 1.876 | 1.895 | 1.844 | 2.072 | | | | 1.910 | 10.17 |
| 1,2-dichloropropane(sim) | 0.438 | 0.513 | 0.406 | 0.413 | 0.521 | 0.445 | 0.407 | 0.413 | 0.399 | 0.453 | | | | 0.441 | 9.98 |
| Bromodichloromethane(sim) | 0.535 | 0.533 | 0.550 | 0.540 | 0.667 | 0.535 | 0.500 | 0.535 | 0.499 | 0.569 | | | | 0.546 | 8.63 |
| Trichloroethene(sim) | 0.509 | 0.424 | 0.414 | 0.396 | 0.544 | 0.471 | 0.434 | 0.443 | 0.421 | 0.476 | | | | 0.453 | 10.23 |
| 1,4-Dioxane(sim) | | | | | 0.211 | 0.188 | 0.173 | 0.174 | 0.156 | 0.177 | | | | 0.180 | 10.15 |
| cis-1,3-Dichloropropene(sim) | 0.481 | 0.448 | 0.421 | 0.425 | 0.539 | 0.475 | 0.452 | 0.466 | 0.449 | | | | | 0.462 | 7.71 |
| 1,1,2-Trichloroethane(sim) | 0.327 | 0.362 | 0.340 | 0.355 | 0.406 | 0.356 | 0.324 | 0.344 | 0.330 | 0.368 | | | | 0.351 | 7.00 |
| Dibromochloromethane(sim) | 0.622 | 0.531 | 0.657 | 0.640 | 0.775 | 0.660 | 0.616 | 0.674 | 0.631 | | | | | 0.645 | 9.90 |
| 1,2-Dibromoethane(EDB)(sim) | 0.634 | 0.542 | 0.465 | 0.521 | 0.600 | 0.574 | 0.520 | 0.547 | 0.529 | 0.599 | | | | 0.553 | 8.89 |
| Tetrachloroethene(sim) qfi | 0.022 | 0.040 | 0.069 | 0.099 | 0.112 | 0.198 | 0.464 | 0.952 | 2.293 | 5.220 | 9.946 | | | Coef R2 | 1.00 |
| Bromoform(sim) | 1.233 | 1.031 | 1.012 | 0.996 | 1.329 | 1.157 | 1.125 | 1.304 | 1.226 | | | | | 1.157 | 10.85 |
| m,p-Xylene(sim) | 3.08 | 2.703 | 2.516 | 2.511 | 2.766 | 2.337 | 2.170 | 2.240 | 2.118 | 2.305 | | | | 2.475 | 12.27 |
| 1,1,2,2-Tetrachloroethane(sim) | 1.964 | 1.570 | 1.439 | 1.427 | 1.864 | 1.652 | 1.555 | 1.615 | 1.551 | 1.591 | | | | 1.623 | 10.49 |
| Benzyl chloride(sim) qfi | 0.012 | 0.017 | 0.028 | 0.050 | 0.111 | 0.169 | 0.462 | 1.088 | 2.477 | | | | | Coef R2 | 1.00 |
| 1,3-Dichlorobenzene(sim) | 2.044 | 1.290 | 1.155 | 1.153 | 1.716 | 1.558 | 1.558 | 1.745 | 1.732 | 1.727 | | | | 1.568 | 18.44 |
| 1,4-Dichlorobenzene(sim) | 1.893 | 1.604 | 1.241 | 1.290 | 1.437 | 1.194 | 1.146 | 1.268 | 1.271 | | | | | 1.372 | 17.42 |
| sec-Butylbenzene(sim) | | | 3.276 | 3.277 | 4.921 | 4.328 | 4.141 | 4.205 | 4.053 | 3.942 | | | | 4.018 | 13.53 |

(#) The maximum %RSD was not met for this compound

Note: m,p-xylene TV is 2 times the TV Listed

(l) linear (q) quadratic (i) inverse conc weight (i2) inverse conc weight squared (f) force through zero

Compounds not using average response (l, li, lfi, li2, lfi2, q, qi, qfi, qi2, qfi2) display concentrations and not response factors

3
AIR ICV RECOVERY

Lab Name: Phoenix Environmental Labs Client: ACT

Lab Code: Phoenix Case No: _____ SAS No: _____ SDG No GCL97021

ICV - Client Id: ICV_CHEM20_0719

| COMPOUND | SPIKE ADDED (ppbv) | | ICV CONCENTRATION (ppbv) | ICV % REC # | QC. LIMITS REC. | |
|-------------------------------|--------------------|--|--------------------------|-------------|-----------------|-----|
| Propylene | 10 | | 9.945 | 99 | 70 | 130 |
| Dichlorodifluoromethane | 10 | | 9.325 | 93 | 70 | 130 |
| Chloromethane | 10 | | 9.528 | 95 | 70 | 130 |
| 1,2-Dichlorotetrafluoroethane | 10 | | 9.907 | 99 | 70 | 130 |
| Vinyl Chloride | 10 | | 9.577 | 96 | 70 | 130 |
| 1,3-Butadiene | 10 | | 9.504 | 95 | 70 | 130 |
| Bromomethane | 10 | | 9.282 | 93 | 70 | 130 |
| Chloroethane | 10 | | 8.693 | 87 | 70 | 130 |
| Ethanol | 10 | | 17.36 | 174 * | 70 | 130 |
| Acetone | 10 | | 9.823 | 98 | 70 | 130 |
| Trichlorofluoromethane | 10 | | 9.903 | 99 | 70 | 130 |
| Isopropylalcohol | 10 | | 10.89 | 109 | 70 | 130 |
| Acrylonitrile | 10 | | 9.576 | 96 | 70 | 130 |
| 1,1-Dichloroethene | 10 | | 9.784 | 98 | 70 | 130 |
| Methylene Chloride | 10 | | 9.416 | 94 | 70 | 130 |
| Carbon Disulfide | 10 | | 9.735 | 97 | 70 | 130 |
| Trichlorotrifluoroethane | 10 | | 9.993 | 100 | 70 | 130 |
| Trans-1,2-Dichloroethene | 10 | | 9.943 | 99 | 70 | 130 |
| 1,1-Dichloroethane | 10 | | 9.898 | 99 | 70 | 130 |
| Methyl tert-butyl ether(MTBE) | 10 | | 9.645 | 96 | 70 | 130 |
| Methyl Ethyl Ketone | 10 | | 10.11 | 101 | 70 | 130 |
| Cis-1,2-Dichloroethene | 10 | | 10.59 | 106 | 70 | 130 |
| Hexane | 10 | | 9.786 | 98 | 70 | 130 |
| Chloroform | 10 | | 9.754 | 98 | 70 | 130 |
| Ethyl acetate | 10 | | 9.963 | 100 | 70 | 130 |
| Tetrahydrofuran | 10 | | 10.38 | 104 | 70 | 130 |
| 1,2-Dichloroethane | 10 | | 9.980 | 100 | 70 | 130 |
| 1,1,1-Trichloroethane | 10 | | 10.02 | 100 | 70 | 130 |
| Benzene | 10 | | 9.902 | 99 | 70 | 130 |
| Carbon Tetrachloride | 10 | | 10.47 | 105 | 70 | 130 |
| Cyclohexane | 10 | | 9.747 | 97 | 70 | 130 |
| 1,2-dichloropropane | 10 | | 9.826 | 98 | 70 | 130 |
| Bromodichloromethane | 10 | | 10.54 | 105 | 70 | 130 |
| Trichloroethene | 10 | | 11.11 | 111 | 70 | 130 |
| 1,4-Dioxane | 10 | | 12.94 | 129 | 70 | 130 |
| Heptane | 10 | | 9.756 | 98 | 70 | 130 |
| cis-1,3-Dichloropropene | 10 | | 10.59 | 106 | 70 | 130 |
| 4-Methyl-2-pentanone(MIBK) | 10 | | 10.46 | 105 | 70 | 130 |
| trans-1,3-Dichloropropene | 10 | | 10.46 | 105 | 70 | 130 |
| 1,1,2-Trichloroethane | 10 | | 10.47 | 105 | 70 | 130 |
| Toluene | 10 | | 10.29 | 103 | 70 | 130 |
| Dibromochloromethane | 10 | | 10.98 | 110 | 70 | 130 |
| 2-Hexanone(MBK) | 10 | | 11.13 | 111 | 70 | 130 |
| 1,2-Dibromoethane(EDB) | 10 | | 10.83 | 108 | 70 | 130 |

Data Path : H:\AIR2022\CHEM20\07JUL\19\
 Data File : 0719_18.D
 Acq On : 20 Jul 2022 12:30 am
 Operator :
 Sample : 10ppb lcs ; AIR
 Misc :
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Jul 21 12:22:00 2022
 Quant Method : H:\AIR2022\CHEM20\METHODS\20_AIR_0719.M
 Quant Title : VOA Standards for 5 point calibration
 QLast Update : Thu Jul 21 12:19:06 2022
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|----------------|----------|--------|---------|----------|
| Internal Standards | | | | | | |
| 1) Bromochloromethane | 7.689 | 130 | 282259 | 10.000 | ng | 0.00 |
| 37) 1,4-Difluorobenzene | 8.840 | 114 | 1066461 | 10.000 | ng | 0.00 |
| 54) Chlorobenzene-d5 | 11.291 | 82 | 534030 | 10.000 | ng | 0.00 |
| 81) Bromochloromethane(sim) | 7.694 | 130 | 304041 | 10.000 | ng | # 0.00 |
| 96) 1,4-Difluorobenzene(sim) | 8.840 | 114 | 1066148 | 10.000 | ng | 0.00 |
| 106) Chlorobenzene-d5(sim) | 11.291 | 82 | 534030 | 10.000 | ng | # 0.00 |
| System Monitoring Compounds | | | | | | |
| 63) % Bromofluorobenzene | 12.111 | 95 | 725998 | 10.037 | ppbv | 0.00 |
| Spiked Amount | 10.000 | Range 70 - 130 | Recovery | = | 100.40% | |
| Target Compounds | | | | | | |
| | | | | | | Qvalue |
| 2) Propylene | 4.200 | 41 | 253260 | 9.945 | ppbv | 95 |
| 3) Dichlorodifluoromethane | 4.275 | 85 | 617143 | 9.325 | ppbv | 98 |
| 4) Chloromethane | 4.437 | 50 | 310608 | 9.528 | ppbv | 100 |
| 5) 1,2-Dichlorotetrafluor... | 4.534 | 85 | 629489 | 9.907 | ppbv | 91 |
| 6) Vinyl Chloride | 4.642 | 62 | 290566 | 9.577 | ppbv | 98 |
| 7) 1,3-Butadiene | 4.771 | 54 | 270106 | 9.504 | ppbv# | 94 |
| 8) Bromomethane | 5.008 | 94 | 228637 | 9.282 | ppbv# | 95 |
| 9) Chloroethane | 5.148 | 64 | 125994 | 8.693 | ppbv | 97 |
| 10) Vinyl Bromide | 5.428 | 106 | 249855 | 9.829 | ppbv# | 94 |
| 11) Ethanol | 5.234 | 45 | 230393 | 17.358 | ppbv | 94 |
| 12) Acetone | 5.612 | 43 | 545983 | 9.823 | ppbv# | 80 |
| 13) Trichlorofluoromethane | 5.752 | 101 | 680083 | 9.903 | ppbv | 100 |
| 14) Isopropylalcohol | 5.795 | 45 | 761716 | 10.886 | ppbv# | 90 |
| 15) Acrylonitrile | 5.946 | 53 | 278452 | 9.576 | ppbv | 91 |
| 16) 1,1-Dichloroethene | 6.202 | 61 | 521137 | 9.784 | ppbv | 96 |
| 17) Methylene Chloride | 6.280 | 49 | 418399 | 9.416 | ppbv# | 81 |
| 18) tert-butyl alcohol | 6.237 | 59 | 721696 | 10.456 | ppbv# | 95 |
| 19) Allyl Chloride | 6.366 | 41 | 446144 | 10.162 | pbv # | 84 |
| 20) Carbon Disulfide | 6.521 | 76 | 654152 | 9.735 | ppbv | 99 |
| 21) Trichlorotrifluoroethane | 6.461 | 101 | 577674 | 9.993 | ppbv | 95 |
| 22) Trans-1,2-Dichloroethene | 6.912 | 61 | 482509 | 9.942 | ppbv | 86 |
| 23) 1,1-Dichloroethane | 7.046 | 63 | 563421 | 9.898 | ppbv | 94 |
| 24) Methyl tert-butyl ethe... | 7.078 | 73 | 651292 | 9.645 | ppbv# | 84 |
| 25) Vinyl Acetate | 7.109 | 43 | 1964384 | 11.007 | ppbv# | 97 |
| 26) Methyl Ethyl Ketone | 7.283 | 43 | 770145 | 10.108 | ppbv# | 92 |
| 27) Cis-1,2-Dichloroethene | 7.584 | 61 | 514407 | 10.587 | ppbv# | 82 |
| 28) Hexane | 7.699 | 57 | 579718 | 9.786 | ppbv# | 83 |
| 29) Chloroform | 7.761 | 83 | 548744 | 9.754 | ppbv | 95 |
| 30) Ethyl acetate | 7.678 | 61 | 126766 | 9.963 | ppbv# | 77 |
| 31) Tetrahydrofuran | 8.022 | 42 | 460428 | 10.378 | ppbv# | 85 |
| 32) 1,2-Dichloroethane | 8.230 | 62 | 426918 | 9.980 | ppbv | 95 |
| 33) 1,1,1-Trichloroethane | 8.386 | 97 | 591354 | 10.017 | ppbv | 95 |
| 34) Benzene | 8.670 | 78 | 786052 | 9.902 | ppbv# | 94 |
| 35) Carbon Tetrachloride | 8.760 | 117 | 639653 | 10.472 | ppbv | 96 |
| 36) Cyclohexane | 8.840 | 84 | 295210 | 9.747 | ppbv# | 69 |
| 38) 1,2-dichloropropane | 9.134 | 63 | 422307 | 9.826 | ppbv# | 84 |
| 39) Bromodichloromethane | 9.247 | 83 | 596399 | 10.545 | ppbv | 98 |
| 40) Trichloroethene | 9.270 | 130 | 476782 | 11.110 | ppbv | 96 |
| 41) 2,2,4-trimethylpentane | 9.270 | 57 | 1863311 | 10.005 | ppbv | 96 |
| 42) 1,4-Dioxane | 9.258 | 88 | 231749 | 12.942 | ppbv# | 73 |
| 43) Methyl methacrylate | 9.326 | 69 | 308769 | 10.575 | ppbv | 99 |
| 44) Heptane | 9.394 | 43 | 758878 | 9.756 | ppbv# | 89 |

Data Path : H:\AIR2022\CHEM20\07JUL\19\
 Data File : 0719_18.D
 Acq On : 20 Jul 2022 12:30 am
 Operator :
 Sample : 10ppb lcs ; AIR
 Misc :
 ALS Vial : 1 Sample Multiplier: 1

Quant Time: Jul 21 12:22:00 2022
 Quant Method : H:\AIR2022\CHEM20\METHODS\20_AIR_0719.M
 Quant Title : VOA Standards for 5 point calibration
 QLast Update : Thu Jul 21 12:19:06 2022
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------------------|--------|------|----------|--------|-------|----------|
| 45) cis-1,3-Dichloropropene | 9.745 | 75 | 489092 | 10.589 | ppbv | 98 |
| 46) 4-Methyl-2-pentanone(M... | 9.745 | 43 | 1012760 | 10.462 | ppbv# | 92 |
| 47) trans-1,3-Dichloropropene | 10.028 | 75 | 446825 | 10.457 | ppbv | 97 |
| 48) 1,1,2-Trichloroethane | 10.141 | 97 | 379915 | 10.469 | ppbv | 97 |
| 49) Toluene | 10.311 | 91 | 1132803 | 10.286 | ppbv | 99 |
| 50) Dibromochloromethane | 10.553 | 129 | 867426 | 10.984 | ppbv | 99 |
| 51) 2-Hexanone(MBK) | 10.413 | 43 | 1053828 | 11.128 | ppbv# | 96 |
| 52) 1,2-Dibromoethane(EDB) | 10.699 | 107 | 641411 | 10.833 | ppbv | 99 |
| 53) Tetrachloroethene | 10.942 | 166 | 622237 | 10.927 | ppbv | 98 |
| 55) 1,1,1,2-Tetrachloroethane | 11.301 | 131 | 491106 | 10.400 | ppbv | 100 |
| 56) Chlorobenzene | 11.312 | 112 | 935978 | 10.475 | ppbv | 95 |
| 57) Ethylbenzene | 11.496 | 91 | 1462258 | 9.840 | ppbv | 99 |
| 58) m,p-Xylene | 11.588 | 91 | 2330160 | 20.676 | ppbv | 98 |
| 59) Bromoform | 11.670 | 173 | 759250 | 12.146 | ppbv | 98 |
| 60) Styrene | 11.793 | 104 | 906266 | 10.974 | ppbv# | 91 |
| 61) 1,1,2,2-Tetrachloroethane | 11.855 | 83 | 826114 | 10.854 | ppbv | 96 |
| 62) o-Xylene | 11.855 | 91 | 1195629 | 10.187 | ppbv | 97 |
| 64) 2-Chlorotoluene | 12.460 | 126 | 405612 | 10.010 | ppbv | 89 |
| 65) Isopropylbenzene | 12.173 | 105 | 1530665 | 10.015 | ppbv | 97 |
| 66) n-Propylbenzene | 12.460 | 120 | 427316 | 10.046 | ppbv | 98 |
| 67) 4-Ethyltoluene | 12.542 | 105 | 1920045 | 11.409 | ppbv | 94 |
| 68) 1,3,5-Trimethylbenzene | 12.583 | 105 | 1411350 | 11.053 | ppbv | 89 |
| 69) 1,2,4-Trimethylbenzene | 12.829 | 105 | 1590219 | 11.725 | ppbv# | 94 |
| 70) tert-butylbenzene | 12.829 | 119 | 1485688 | 9.849 | ppbv | 96 |
| 71) Benzyl chloride | 12.932 | 91 | 996589 | 14.203 | ppbv# | 84 |
| 72) 1,3-Dichlorobenzene | 12.952 | 146 | 1155391 | 13.550 | ppbv | 100 |
| 73) 1,4-Dichlorobenzene | 12.983 | 146 | 856240 | 13.234 | ppbv | 98 |
| 74) sec-Butylbenzene | 12.993 | 105 | 1860495 | 9.589 | ppbv | 98 |
| 75) 4-Isopropyltoluene | 13.075 | 119 | 1769710 | 9.525 | ppbv | 96 |
| 76) 1,2-Dichlorobenzene | 13.209 | 146 | 1051051 | 13.744 | ppbv | 99 |
| 77) n-Butylbenzene | 13.342 | 91 | 1418536 | 9.894 | ppbv | 99 |
| 78) 1,2,4-Trichlorobenzene | 14.399 | 180 | 585884 | 16.714 | ppbv | 99 |
| 79) Naphthalene | 14.492 | 128 | 1440545 | 14.775 | ppbv# | 100 |
| 80) Hexachlorobutadiene | 14.728 | 225 | 886175 | 13.415 | ppbv | 99 |
| 82] 1,2-Dichlorotetrafluor... | 4.534 | 85 | 627047 | 9.011 | ppbv | 91 |
| 83] Vinyl Chloride(sim) | 4.647 | 62 | 311149 | 9.483 | ppbv | 97 |
| 84] Bromomethane(sim) | 5.008 | 94 | 228637 | 8.027 | ppbv# | 95 |
| 85] Trichlorofluoromethane... | 5.757 | 101 | 716060 | 9.872 | ppbv# | 99 |
| 86] 1,2-Dichloroethane(sim) | 8.230 | 62 | 427391 | 9.770 | ppbv | 95 |
| 87] 1,1,1-Trichloroethane(... | 8.392 | 97 | 632124 | 10.201 | ppbv# | 95 |
| 88] Benzene(sim) | 8.670 | 78 | 786498 | 8.461 | ppbv# | 94 |
| 89] Carbon Tetrachloride(sim) | 8.766 | 117 | 672758 | 10.879 | ppbv | 97 |
| 90] 1,1-Dichloroethene(sim) | 6.202 | 61 | 521137 | 9.388 | ppbv | 96 |
| 91] Trichlorotrifluoroetha... | 6.458 | 101 | 607603 | 9.886 | ppbv# | 98 |
| 92] Trans-1,2-Dichloroethe... | 6.912 | 61 | 482509 | 9.697 | ppbv | 86 |
| 93] 1,1-Dichloroethane(sim) | 7.052 | 63 | 603076 | 9.918 | ppbv# | 96 |
| 94] Cis-1,2-Dichloroethene... | 7.584 | 61 | 514407 | 10.745 | ppbv# | 82 |
| 95] Chloroform(sim) | 7.767 | 83 | 580265 | 9.991 | ppbv | 95 |
| 97] 1,2-dichloropropane(sim) | 9.140 | 63 | 454381 | 9.669 | ppbv# | 85 |
| 98] Bromodichloromethane(sim) | 9.247 | 83 | 596399 | 10.240 | ppbv | 98 |
| 99] Trichloroethene(sim) | 9.275 | 130 | 506810 | 10.489 | ppbv | 96 |
| 100] 1,4-Dioxane(sim) | 9.258 | 88 | 231749 | 12.089 | ppbv# | 73 |
| 101] cis-1,3-Dichloropropen... | 9.751 | 75 | 526872 | 10.702 | ppbv | 98 |
| 102] 1,1,2-Trichloroethane(... | 10.141 | 97 | 379915 | 10.147 | ppbv | 97 |
| 103] Dibromochloromethane(sim) | 10.559 | 129 | 813315 | 11.823 | ppbv | 100 |

Data Path : H:\AIR2022\CHEM20\07JUL\19\
 Data File : 0719_18.D
 Acq On : 20 Jul 2022 12:30 am
 Operator :
 Sample : 10ppb lcs ; AIR
 Misc :
 ALS Vial : 1 Sample Multiplier: 1

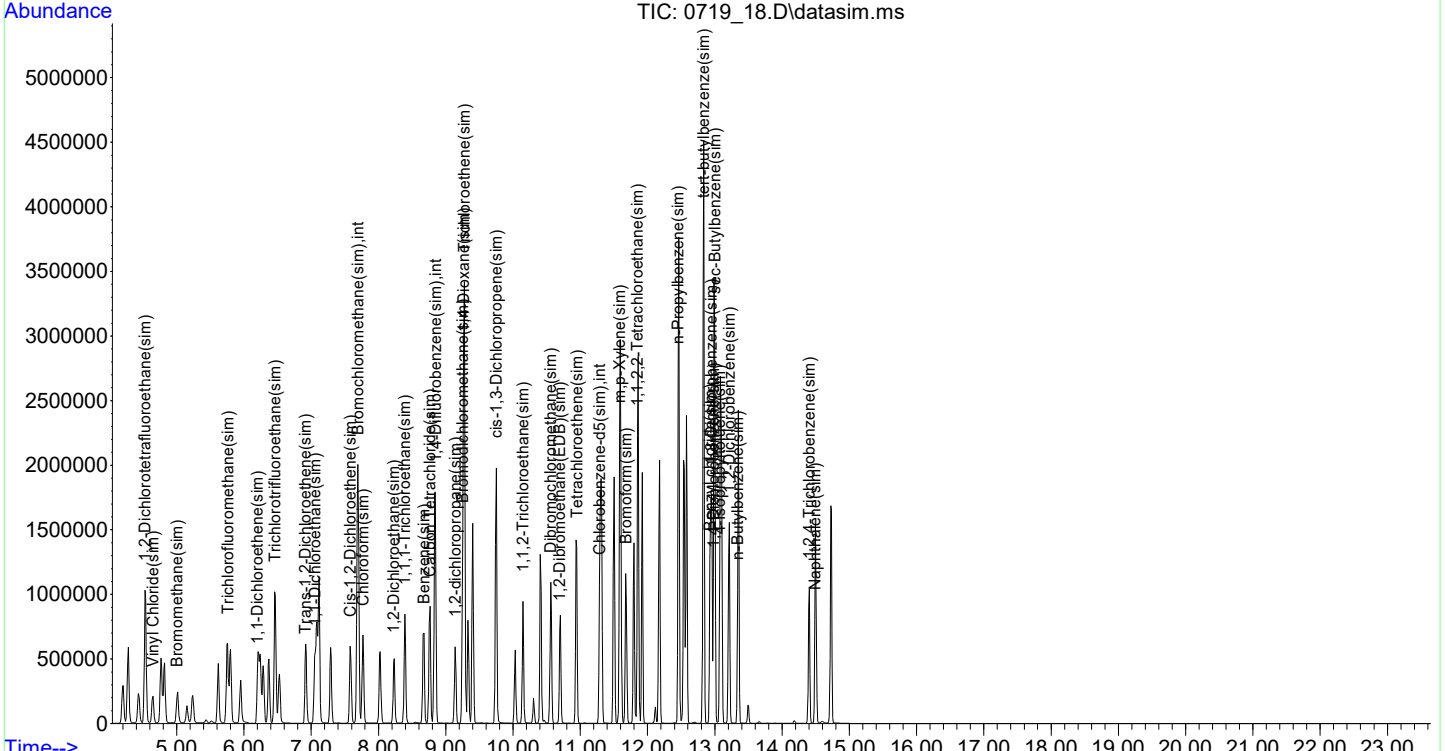
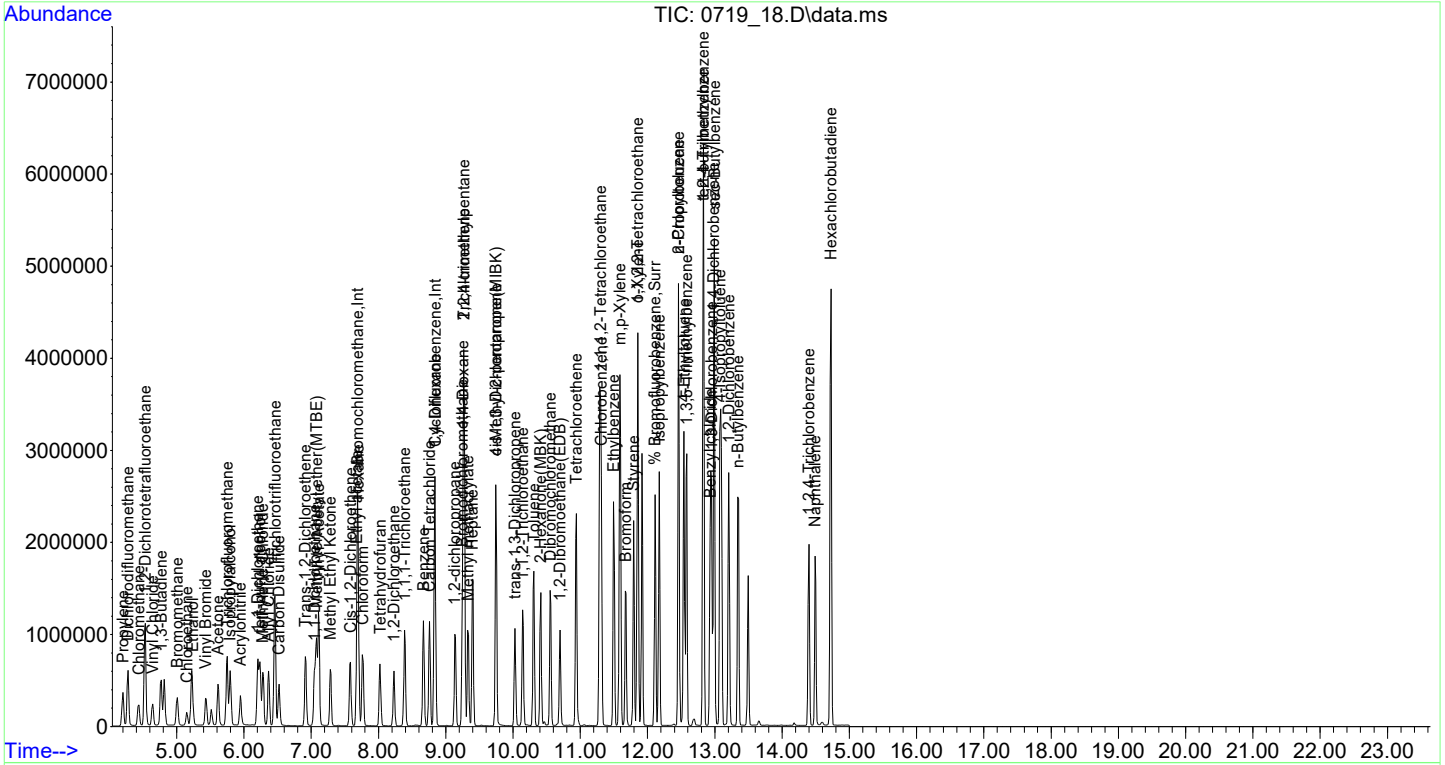
Quant Time: Jul 21 12:22:00 2022
 Quant Method : H:\AIR2022\CHEM20\METHODS\20_AIR_0719.M
 Quant Title : VOA Standards for 5 point calibration
 QLast Update : Thu Jul 21 12:19:06 2022
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------------------|--------|------|----------|-------------|-------|----------|
| 104] 1,2-Dibromoethane(EDB)... | 10.699 | 107 | 641411 | 10.877 | ppbv | 99 |
| 105] Tetrachloroethene(sim) | 10.947 | 166 | 686788 | 10.853 | ppbv | 99 |
| 107] Bromoform(sim) | 11.676 | 173 | 829597 | 13.427 | ppbv | 99 |
| 108] m,p-Xylene(sim) | 11.588 | 91 | 2336280 | 17.678 | ppbv | 98 |
| 109] 1,1,2,2-Tetrachloroeth... | 11.850 | 83 | 862263 | 9.949 | ppbv | 97 |
| 110] n-Propylbenzene(sim) | 12.460 | 120 | 427316 | 9.048 | ppbv | 98 |
| 111] tert-butylbenzene(sim) | 12.835 | 119 | 1596164 | 9.186 | ppbv# | 95 |
| 112] Benzyl chloride(sim) | 12.932 | 91 | 996589 | 10.437 | ppbv | 98 |
| 113] 1,3-Dichlorobenzene(sim) | 12.948 | 146 | 1138811 | 13.603 | ppbv | 99 |
| 114] 1,4-Dichlorobenzene(sim) | 12.983 | 146 | 856240 | 11.690 | ppbv | 98 |
| 115] sec-Butylbenzene(sim) | 12.999 | 105 | 1939945 | 9.041 | ppbv | 97 |
| 116] 4-Isopropyltoluene(sim) | 13.075 | 119 | 1770660 | 8.881 | ppbv | 96 |
| 117] 1,2-Dichlorobenzene(sim) | 13.214 | 146 | 1148034 | 13.871 | ppbv | 99 |
| 118] n-Butylbenzene(sim) | 13.342 | 91 | 1418536 | 9.555 | ppbv | 99 |
| 119] 1,2,4-Trichlorobenzene... | 14.405 | 180 | 654884 | 16.058 | ppbv | 99 |
| 120] Naphthalene(sim) | 14.492 | 128 | 1440545 | 4.410 | ppbv | 99 |
| 121] Hexachlorobutadiene(sim) | 14.733 | 225 | 1043174 | Outside Cal | | 100 |

(#)out of range (m)manual integration reviewed by analyst (+)signals summed

Data Path : H:\AIR2022\CHEM20\07JUL\19\
 Data File : 0719_18.D
 Acq On : 20 Jul 2022 12:30 am
 Operator :
 Sample : 10ppb lcs ; AIR
 Misc :
 ALS Vial : 1 Sample Multiplier: 1

Quant Time : Jul 21 12:22:00 2022
 Quant Method : H:\AIR2022\CHEM20\METHODS\20_AIR_0719.M
 Quant Title : VOA Standards for 5 point calibration
 QLast Update : Thu Jul 21 12:19:06 2022
 Response via : Initial Calibration



ATTACHMENT D

ANNOTATED SUMMARY FORMS

1
AIR ANALYSIS DATA SHEET

CLIENT ID

IA-4

| | | | |
|--------------|----------|------------------|-------------------|
| Client: | ACT | Lab: | Phoenix Env. Labs |
| SDG No.: | GCL97021 | Lab Sample ID: | CL97021 |
| Canister: | 28560 | Lab File ID: | 0803_10.D |
| Instrument: | CHEM20 | Column: | RTX-1 60M |
| Purge Volume | 200 (cc) | Date Received: | 08/03/22 |
| Matrix: | AIR | Date Analyzed: | 08/03/22 |
| | | Dilution Factor: | 1 |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|-------------|-------------------------------|-------|----|-------|-------|---|
| 115-07-1 | Propylene | 0.581 | U | 0.581 | 0.581 | r |
| 75-71-8 | Dichlorodifluoromethane | 0.548 | | 0.202 | 0.202 | r |
| 74-87-3 | Chloromethane | 0.596 | J+ | 0.485 | 0.485 | r |
| 75-01-4 | Vinyl Chloride | 1.82 | | 0.078 | 0.078 | r |
| 106-99-0 | 1,3-Butadiene | 0.452 | U | 0.452 | 0.452 | r |
| 75-00-3 | Chloroethane | 0.379 | U | 0.379 | 0.379 | r |
| 64-17-5 | Ethanol | 62.8 | J | 0.531 | 0.531 | r |
| 67-64-1 | Acetone | 31.0 | J+ | 0.421 | 0.421 | r |
| 75-69-4 | Trichlorofluoromethane | 0.265 | | 0.178 | 0.178 | r |
| 67-63-0 | Isopropylalcohol | 4.48 | | 0.407 | 0.407 | r |
| 107-13-1 | Acrylonitrile | 0.461 | U | 0.461 | 0.461 | r |
| 75-09-2 | Methylene Chloride | 0.863 | U | 0.863 | 0.863 | r |
| 75-15-0 | Carbon Disulfide | 0.321 | U | 0.321 | 0.321 | r |
| 1634-04-4 | Methyl tert-butyl ether(MTBE) | 0.278 | U | 0.278 | 0.278 | r |
| 78-93-3 | Methyl Ethyl Ketone | 1.70 | | 0.339 | 0.339 | r |
| 156-59-2 | Cis-1,2-Dichloroethene | 17.2 | | 0.051 | 0.051 | r |
| 110-54-3 | Hexane | 5.37 | | 0.284 | 0.284 | r |
| 141-78-6 | Ethyl acetate | 0.278 | U | 0.278 | 0.278 | r |
| 109-99-9 | Tetrahydrofuran | 6.13 | | 0.339 | 0.339 | r |
| 71-43-2 | Benzene | 1.28 | | 0.313 | 0.313 | r |
| 110-82-7 | Cyclohexane | | U | 1.81 | 1.81 | r |
| 79-01-6 | Trichloroethene | 2.30 | | 0.037 | 0.037 | r |
| 142-82-5 | Heptane | 1.64 | | 0.244 | 0.244 | r |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | 0.244 | U | 0.244 | 0.244 | r |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.221 | U | 0.221 | 0.221 | r |
| 108-88-3 | Toluene | 7.16 | | 0.266 | 0.266 | r |
| 591-78-6 | 2-Hexanone(MBK) | 0.244 | U | 0.244 | 0.244 | r |
| 127-18-4 | Tetrachloroethene | 3.28 | | 0.037 | 0.037 | r |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.146 | U | 0.146 | 0.146 | r |
| 108-90-7 | Chlorobenzene | 0.217 | U | 0.217 | 0.217 | r |
| 100-41-4 | Ethylbenzene | 0.928 | | 0.230 | 0.230 | r |
| 179601-23-1 | m,p-Xylene | 3.21 | | 0.230 | 0.230 | r |
| 100-42-5 | Styrene | 0.235 | U | 0.235 | 0.235 | r |
| 95-47-6 | o-Xylene | 1.23 | | 0.230 | 0.230 | r |
| 98-82-8 | Isopropylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 622-96-8 | 4-Ethyltoluene | 0.972 | | 0.204 | 0.204 | r |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

1
AIR ANALYSIS DATA SHEET

CLIENT ID

IA-7

| | | | |
|----------------|----------|------------------|-------------------|
| Client: | ACT | Lab: | Phoenix Env. Labs |
| SDG No.: | GCL97021 | Lab Sample ID: | CL97022 |
| Canister: | 23330 | Lab File ID: | 0803_11.D |
| Instrument: | CHEM20 | Column: | RTX-1 60M |
| Date Received: | 08/03/22 | | |
| Purge Volume | 200 (cc) | Date Analyzed: | 08/03/22 |
| Matrix: | AIR | Dilution Factor: | 1 |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|-------------|-------------------------------|-------|----|-------|-------|---|
| 115-07-1 | Propylene | 1.01 | | 0.581 | 0.581 | r |
| 75-71-8 | Dichlorodifluoromethane | 0.502 | | 0.202 | 0.202 | r |
| 74-87-3 | Chloromethane | 0.630 | J+ | 0.485 | 0.485 | r |
| 75-01-4 | Vinyl Chloride | 1.90 | | 0.078 | 0.078 | r |
| 106-99-0 | 1,3-Butadiene | 0.452 | U | 0.452 | 0.452 | r |
| 75-00-3 | Chloroethane | 0.379 | U | 0.379 | 0.379 | r |
| 64-17-5 | Ethanol | 76.1 | J | 0.531 | 0.531 | r |
| 67-64-1 | Acetone | 44.7 | J+ | 0.421 | 0.421 | r |
| 75-69-4 | Trichlorofluoromethane | 0.289 | | 0.178 | 0.178 | r |
| 67-63-0 | Isopropylalcohol | 4.24 | | 0.407 | 0.407 | r |
| 107-13-1 | Acrylonitrile | 0.461 | U | 0.461 | 0.461 | r |
| 75-09-2 | Methylene Chloride | 0.863 | U | 0.863 | 0.863 | r |
| 75-15-0 | Carbon Disulfide | 0.321 | U | 0.321 | 0.321 | r |
| 1634-04-4 | Methyl tert-butyl ether(MTBE) | 0.278 | U | 0.278 | 0.278 | r |
| 78-93-3 | Methyl Ethyl Ketone | 1.17 | | 0.339 | 0.339 | r |
| 156-59-2 | Cis-1,2-Dichloroethene | 16.0 | | 0.051 | 0.051 | r |
| 110-54-3 | Hexane | 6.22 | | 0.284 | 0.284 | r |
| 141-78-6 | Ethyl acetate | 0.278 | U | 0.278 | 0.278 | r |
| 109-99-9 | Tetrahydrofuran | 2.57 | | 0.339 | 0.339 | r |
| 71-43-2 | Benzene | 1.72 | | 0.313 | 0.313 | r |
| 110-82-7 | Cyclohexane | | U | 2.16 | 2.16 | r |
| 79-01-6 | Trichloroethene | 1.90 | | 0.037 | 0.037 | r |
| 142-82-5 | Heptane | 1.99 | | 0.244 | 0.244 | r |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | 0.244 | U | 0.244 | 0.244 | r |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.221 | U | 0.221 | 0.221 | r |
| 108-88-3 | Toluene | 7.10 | | 0.266 | 0.266 | r |
| 591-78-6 | 2-Hexanone(MBK) | 0.244 | U | 0.244 | 0.244 | r |
| 127-18-4 | Tetrachloroethene | 2.80 | | 0.037 | 0.037 | r |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.146 | U | 0.146 | 0.146 | r |
| 108-90-7 | Chlorobenzene | 0.217 | U | 0.217 | 0.217 | r |
| 100-41-4 | Ethylbenzene | 0.995 | | 0.230 | 0.230 | r |
| 179601-23-1 | m,p-Xylene | 3.54 | | 0.230 | 0.230 | r |
| 100-42-5 | Styrene | 0.235 | U | 0.235 | 0.235 | r |
| 95-47-6 | o-Xylene | 1.33 | | 0.230 | 0.230 | r |
| 98-82-8 | Isopropylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 622-96-8 | 4-Ethyltoluene | 1.03 | | 0.204 | 0.204 | r |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

December 8, 2025

Mr. Jason Stewart
Advanced Cleanup Technologies
228 Park Ave S PMB 34864
New York, New York 10003

Re: Data Usability Summary Report – Phoenix Environmental Laboratories, Inc. – GCM85136 V1

Dear Mr. Stewart:

The evaluation of analytical data by Phoenix Environmental Laboratories for project 9268-BXNY, which were reported in a single data package under Job No. GCM85136 V1 has been completed. The following samples were reported.

SS-12 IA-11 OA-2

Analysis was performed in accordance with EPA Method TO-15 (volatile organics). The review was performed to the extent possible, in accordance with the analytical method, and “DER-10/ Technical Guidance for Site Investigation and Remediation”. Professional judgment is applied as necessary and appropriate. National Functional Guidelines for Organic Data Review was consulted as needed. Qualifiers consistent with those defined by EPA Region 2 are applied as necessary and appropriate.

Below is the Data Usability Summary Report (DUSR) associated with these samples.

| Data Usability Summary Report | |
|---|--------------------------------|
| 1. Is the data package complete as defined under the requirements for the most current DEC ASP Category B or USEPA CLP data deliverables? | No – see Documentation Section |
| 2. Have all holding times been met? | Yes |
| 3. Do all the QC data; blanks, instrument tunings, calibration standards, calibration verifications, surrogate recoveries, spike recoveries, replicate analyses, laboratory controls and sample data fall within the protocol required limits and specifications? | No -see following sections |
| 4. Have all of the data been generated using established and agreed upon analytical protocols? | Yes |
| 5. Does an evaluation of the raw data confirm the results provided in the data summary sheet and the quality control verification forms? | Yes |
| 6. Have the correct data qualifiers been used and are they consistent with the most current DEC ASP? | Yes |
| 7. Have any quality control (QC) exceedances been specifically noted in the DUSR and have the corresponding QC summary sheet from the data package been attached to the DUSR? | Yes |

Overall Evaluation

Based on the data review effort, results are usable, with the following qualifications. For samples that are qualified as estimated (J-, UJ), detected results may be biased low. False negatives may exist in non-detect results. Sample results that are qualified as estimated (J+) may be biased high. For samples that are qualified as estimated with any combination of (J), (J-) and/or (J+), the (J) qualifier takes precedence and is applied to the sample result. It is not possible to determine the direction of the bias and the overall effect on the result.

- The results for all target analytes in all samples are qualified as estimated (J, UJ) because there is no evidence that the samples were properly preserved .
- The result for 1,2,4-trichlorobenzene in SS-12 is qualified as estimated (UJ) due to low response in the continuing calibration verification (CCV) standard.
- The results for acetone in IA-11 and OA-2 are qualified as estimated (J+) due to high response in the CCV.
- The results for isopropyl alcohol and 1,4-dioxane in IA-11 and OA-2 are qualified as estimated (J-, UJ) due to low recovery in the laboratory control sample (LCS).
- The results for ethanol in IA-11 and acetone in SS-12 (undiluted analysis) are qualified as estimated (J) because the reported concentration is outside the calibration range.

Qualifier definitions are provided in Attachment A. A copy of the chain of custody record is provided in Attachment B. Pages from the data package illustrating the exceedances and issues described in this validation report are included in Attachment C. Annotated Summary Forms are included in Attachment D detailing qualifications resulting from the data review effort.

The following components were reviewed, where applicable:

- Chain of Custody
- Receiving conditions
- Holding times
- Preservation
- Analyte lists
- Reporting limits
- Requested methods
- Units, and
- Sample related quality control data:
 - Method, instrument blanks
 - Clean canister certification
 - Field blanks

- Surrogate recoveries
- LCS recoveries
- Internal standard area response
- Duplicates
- Analyte Identification
- Instrument related quality control data:
 - Instrument tunes
 - Calibration summaries

The following sections of the report detail only quality control exceedances that impacted results. Where a quality control item exceeded control limits but there is no impact to the samples results, these are not detailed in the report.

Documentation: A completeness review of the data package was performed, and the data package was determined to be a complete Category B data package, with the following exceptions:

- The data package does not include summary forms and raw data for the second source initial calibration verification (ICV) standard. The laboratory was contacted and provided the ICV data and is included in attachment C.
- Calibration data for the clean canister certification analysis is not included in the data package.

The following documentation issues were observed during the review:

- Results for ethanol, acetone, isopropyl alcohol, and hexane in all samples are flagged 'S' which the data deliverable identifies this flag as 'This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level.' Where the sample detections are greater than five times the reporting limit, the validator removed the 'S' qualifier.

Holding Times, Preservation, Sample Integrity:

A copy of the applicable chain of custody (COC) record was included in the data package, documenting sample collection date of November 9, 2022. The samples were received on November 14, 2022. Samples were analyzed within the method specified hold time. However, the data package does contain sample receipt information regarding condition of canisters upon receipt. Since there is no evidence that the samples were properly preserved (pressure inside the canister maintained within +/- 5 psi from sampling to check in the laboratory or analysis), the results for all target analytes in all samples are qualified as estimated (J, UJ).

A. Volatile Organics

1. Calibration

Two initial calibrations (IC) were performed in support of the sample analysis. All relative response factors (RRFs) and relative standard deviations (RSDs) or correlation coefficients are acceptable. A second source ICV standard was analyzed after each IC, and all percent differences are acceptable (70-130%R). Continuing calibration verification (CCV) standard were analyzed at the appropriate frequency and are acceptable (%D<30), with the exceptions noted below

| Analyte | %D | Associated Sample | Qualifier Applied |
|----------------------------------|-------|-------------------|-------------------|
| <i>CCV CHEM20 11/14/22 12:47</i> | | | |
| 1,2,4-Trichlorobenzene (SIM) | 34.4 | SS-12 | UJ |
| <i>CCV CHEM39 11/14/22 14:40</i> | | | |
| Acetone | -35.5 | IA-11 OA-2 | J+ |

The percent difference for 1,2,4-trichlorobenzene (SIM) represents a decrease in instrument sensitivity. The result for 1,2,4-trichlorobenzene in SS-12 is qualified as estimated (UJ) due to low response in the CCV.

The percent difference for acetone represents an increase in instrument sensitivity. The results for acetone in IA-11 and OA-2 are qualified as estimated (J+) due to high response in the CCV.

2. Laboratory Control Sample (LCS)

Two LCSs were prepared and analyzed with the sample. The LCS is evaluated using control limit of 70-130%R. All recoveries are acceptable, with the following exceptions:

| Analyte | LCS %R | Affected Sample | Qualifier Applied |
|--------------------|--------|-----------------|-------------------|
| <i>CM85137 LCS</i> | | | |
| Isopropyl alcohol | 63 | IA-11 | J-, UJ |
| 1,4-Dioxane (SIM)* | 64* | OA-2 | |

*The summary form provided by the laboratory reports LCS recoveries based on the full scan analysis. The validator reviewed the raw data to determine the recovery in the SIM analysis and this recovery is presented above.

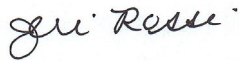
The results for isopropyl alcohol and 1,4-dioxane (SIM) in IA-11 and OA-2 are qualified as estimated (J-, UJ) due to low recovery in the LCS.

3. Compound Quantitation

The result for ethanol in IA-11 and result for acetone in SS-12 (undiluted analysis) is qualified as estimated (J) because the reported concentration is outside the calibration range. Results from the diluted analysis.

No other sample results are qualified. Please feel free to contact me at (908) 370-3431 or richjერიrossi513@gmail.com if you have any questions regarding this data package review report or need further information.

Sincerely,

A handwritten signature in cursive script that reads "Jeri Rossi".

Jeri L Rossi, CEAC

Environmental Consulting Chemist

ATTACHMENT A

Qualifier Definitions

EPA Qualifier Definitions

- U The analyte was analyzed for but was not detected above the level of the reported sample quantitation limit.
- J The result is an estimated quantity. The associated numerical value is the approximate concentration of the analyte in the sample.
- J+ The result is an estimated quantity, but the result may be biased high.
- J- The result is an estimated quantity, but the result may be biased low.
- NJ The analyte has been “tentatively identified” or “presumptively” as present and the associated numerical value is the estimated concentration in the sample.
- UJ The analyte was analyzed for but was not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
- R The data are unusable. The sample results are rejected due to serious deficiencies in meeting QC criteria. The analyte may or may not be present in the sample.

ATTACHMENT B

CHAIN OF CUSTODY (COC)



11720 Middlebrook Rd. Suite 110, Madison, NJ 07952
 Telephone: 973-645-1100 Fax: 973-645-0071

CHAIN OF CUSTODY RECORD

AIR ANALYSES

860-645-1102

email: greg@phoenixlabs.com

P.O. #

Page of

Data Delivery:

Fax #: _____
 Email: _____
 Phone #: _____

| | | | |
|---|-----------------------------------|---|--|
| Report to: <u>Jason Stewart</u> | Project Name: <u>9628-BXNY</u> | Data Format: (Circle) Equis <u>Excel</u> Other: | Ambient/Indoor Air Soil Gas Grab (G) Composite (C) TO-15 APH |
| Customer: <u>Advanced Cleanup Technologies, Inc.</u> | Invoice to: <u>Karen Freedman</u> | Requested Deliverable: RCP <u>ASP CAT B</u> | |
| Address: <u>228 Park Ave S</u> <u>PMB 34864 NY, NY 10003</u> | Sampled by: <u>Tim Young</u> | Quote Number: | |

| Phoenix ID # | Client Sample ID | Canister ID # | Canister Size (L) | Outgoing Canister Pressure ("Hg) | Incoming Canister Pressure ("Hg) | Flow Regulator ID # | Flow Controller Setting (mL/min) | Sampling Start Time | Sampling End Time | Sample Start Date | Canister Pressure at Start ("Hg) | Canister Pressure at End ("Hg) | MATRIX | ANALYSES |
|-------------------------------|------------------|---------------|-------------------|----------------------------------|----------------------------------|---------------------|----------------------------------|---------------------|-------------------|-------------------|----------------------------------|--------------------------------|----------|----------|
| THIS SECTION FOR LAB USE ONLY | | | | | | | | | | | | | | |
| <u>85136</u> | <u>SS-12</u> | <u>11287</u> | <u>6.0L</u> | <u>-30</u> | <u>-9</u> | <u>5620</u> | <u>22.8</u> | <u>0945</u> | <u>1240</u> | <u>11/9</u> | <u>-30</u> | <u>-8</u> | <u>X</u> | <u>✓</u> |
| <u>85137</u> | <u>IA-11</u> | <u>28614</u> | <u>6.0L</u> | <u>-30</u> | <u>-8</u> | <u>5618</u> | <u>3.6</u> | <u>1420</u> | <u>0915</u> | <u>11/9/11</u> | <u>-30</u> | <u>-8</u> | <u>X</u> | <u>✓</u> |
| <u>85138</u> | <u>OA-2</u> | <u>11257</u> | <u>6.0L</u> | <u>-30</u> | <u>-5</u> | <u>3248</u> | <u>3.7</u> | <u>1430</u> | <u>0917</u> | <u>11/9/11</u> | <u>-30</u> | <u>-8</u> | <u>X</u> | <u>✓</u> |

| | | | | |
|-------------------------------------|---------------------------------|-----------------------|--------------------|---|
| Relinquished by: <u>[Signature]</u> | Accepted by: <u>[Signature]</u> | Date: <u>11.14.22</u> | Time: <u>10:43</u> | I attest that all media released by Phoenix Environmental Laboratories, Inc. have been received in good working condition and agree to the terms and conditions as listed on the back of this document. |
| | | Date: <u>11.14.22</u> | Time: <u>17:15</u> | |

| | | | | | | | |
|---|---|---|--|--|----------------------------|--|---|
| State Where Samples Collected: <u>Bronx, NY</u> | Turnaround Time: 1 Day <input type="checkbox"/> 2 Day <input type="checkbox"/> 3 Day <input type="checkbox"/> 4 Day <input type="checkbox"/> 5 Day <input checked="" type="checkbox"/> | Requested Criteria: (Please Circle) CT: TAC I/C TAC RES SVVC I/C SVVC RES GWV I/C GWV CES | MA: Indoor Air: Residential Ind/Commercial Soil Gas: Residential Ind/Commercial | NI: Indoor Air: Residential Ind/Commercial Soil Gas: Residential Ind/Commercial | NY: <u>Vapor Intrusion</u> | PA: Indoor Air: Residential Non-residential | VT: Indoor Air: Residential Industrial Sub-slab Residential Industrial |
|---|---|---|--|--|----------------------------|--|---|

SPECIAL INSTRUCTIONS, OC REQUIREMENTS, REGULATORY INFORMATION:

01/09/2024

Phoenix Environmental Laboratories, Inc.

Page 24 of 323

ATTACHMENT C

**SELECTED PAGES FROM DATA PACKAGE –
QC EXCEEDANCES AND VALIDATION ISSUES**

3
AIR LCS RECOVERY

Lab Name: Phoenix Environmental Labs Client: ACT

Lab Code: Phoenix Case No: _____ SAS No: _____ SDG No GCM85136

LCS - Client Id: CM85137 LCS

| COMPOUND | SPIKE ADDED (ppbv) | | LCS CONCENTRATION (ppbv) | LCS % REC # | QC. LIMITS REC. | |
|-------------------------------|--------------------|--|--------------------------|-------------|-----------------|-----|
| Propylene | 10 | | 11.85 | 119 | 70 | 130 |
| Dichlorodifluoromethane | 10 | | 10.34 | 103 | 70 | 130 |
| Chloromethane | 10 | | 11.04 | 110 | 70 | 130 |
| 1,2-Dichlorotetrafluoroethane | 10 | | 10.77 | 108 | 70 | 130 |
| Vinyl Chloride | 10 | | 10.62 | 106 | 70 | 130 |
| 1,3-Butadiene | 10 | | 10.82 | 108 | 70 | 130 |
| Bromomethane | 10 | | 9.582 | 96 | 70 | 130 |
| Chloroethane | 10 | | 10.36 | 104 | 70 | 130 |
| Ethanol | 10 | | 7.239 | 72 | 70 | 130 |
| Acetone | 10 | | 10.72 | 107 | 70 | 130 |
| Trichlorofluoromethane | 10 | | 9.728 | 97 | 70 | 130 |
| Isopropylalcohol | 10 | | 6.293 | 63 * | 70 | 130 |
| Acrylonitrile | 10 | | 10.12 | 101 | 70 | 130 |
| 1,1-Dichloroethene | 10 | | 9.973 | 100 | 70 | 130 |
| Methylene Chloride | 10 | | 10.63 | 106 | 70 | 130 |
| Carbon Disulfide | 10 | | 10.12 | 101 | 70 | 130 |
| Trichlorotrifluoroethane | 10 | | 10.09 | 101 | 70 | 130 |
| Trans-1,2-Dichloroethene | 10 | | 14.69 | 147 * | 70 | 130 |
| 1,1-Dichloroethane | 10 | | 10.54 | 105 | 70 | 130 |
| Methyl tert-butyl ether(MTBE) | 10 | | 9.849 | 98 | 70 | 130 |
| Methyl Ethyl Ketone | 10 | | 10.54 | 105 | 70 | 130 |
| Cis-1,2-Dichloroethene | 10 | | 10.47 | 105 | 70 | 130 |
| Hexane | 10 | | 11.08 | 111 | 70 | 130 |
| Chloroform | 10 | | 10.19 | 102 | 70 | 130 |
| Ethyl acetate | 10 | | 9.101 | 91 | 70 | 130 |
| Tetrahydrofuran | 10 | | 10.70 | 107 | 70 | 130 |
| 1,2-Dichloroethane | 10 | | 9.959 | 100 | 70 | 130 |
| 1,1,1-Trichloroethane | 10 | | 9.772 | 98 | 70 | 130 |
| Benzene | 10 | | 10.15 | 102 | 70 | 130 |
| Carbon Tetrachloride | 10 | | 9.987 | 100 | 70 | 130 |
| Cyclohexane | 10 | | 9.945 | 99 | 70 | 130 |
| 1,2-dichloropropane | 10 | | 11.02 | 110 | 70 | 130 |
| Bromodichloromethane | 10 | | 10.61 | 106 | 70 | 130 |
| Trichloroethene | 10 | | 10.55 | 106 | 70 | 130 |
| 1,4-Dioxane | 10 | | 6.354 | 64 * | 70 | 130 |
| Heptane | 10 | | 12.03 | 120 | 70 | 130 |
| cis-1,3-Dichloropropene | 10 | | 11.15 | 112 | 70 | 130 |
| 4-Methyl-2-pentanone(MIBK) | 10 | | 11.20 | 112 | 70 | 130 |
| trans-1,3-Dichloropropene | 10 | | 10.94 | 109 | 70 | 130 |
| 1,1,2-Trichloroethane | 10 | | 10.63 | 106 | 70 | 130 |
| Toluene | 10 | | 10.79 | 108 | 70 | 130 |
| Dibromochloromethane | 10 | | 11.19 | 112 | 70 | 130 |
| 2-Hexanone(MBK) | 10 | | 10.68 | 107 | 70 | 130 |
| 1,2-Dibromoethane(EDB) | 10 | | 11.02 | 110 | 70 | 130 |

FORM III AIR

7B
AIR CONTINUING CALIBRATION CHECK

Lab Name: Phoenix Environmental Labs Client: _____
 Lab Code: Phoenix Case No.: _____ SAS No.: _____ SDG No.: GCM85136
 Instrument: CHEM20 Calibration Date: 11/14/22 Time: 12:47
 Lab File Id: 1114_02.D Init. Calib. Date(s): 10/10/22 10/10/22
 Heated Purge (Y/N): Y Init. Calib. Times: 11:30 18:12
 GC Column: RTX-1 60M Method File: 20_AIR_1010.M

| COMPOUND | RRF | RRF1 | RRF MIN | %D | % D LIMITS |
|--------------------------------|-------|-------|---------|--------|------------|
| 1,2-Dichloroethane(sim) | 1.535 | 1.840 | | -19.9 | 30 |
| 1,1,1-Trichloroethane(sim) | 2.249 | 2.524 | | -12.2 | 30 |
| Benzene(sim) | 2.853 | 2.357 | | 17.4 | 30 |
| Carbon Tetrachloride(sim) | 2.338 | 2.746 | | -17.5 | 30 |
| 1,1-Dichloroethene(sim) | 1.753 | 1.896 | | -8.2 | 30 |
| Trichlorotrifluoroethane(sim) | 2.090 | 1.854 | | 11.3 | 30 |
| Trans-1,2-Dichloroethene(sim) | 1.726 | 1.789 | | -3.7 | 30 |
| 1,1-Dichloroethane(sim) | 2.069 | 2.243 | | -8.4 | 30 |
| Cis-1,2-Dichloroethene(sim) | 1.602 | 1.723 | | -7.6 | 30 |
| Chloroform(sim) | 2.058 | 2.193 | | -6.6 | 30 |
| 1,2-dichloropropane(sim) | 0.453 | 0.420 | | 7.3 | 30 |
| Bromodichloromethane(sim) | 0.630 | 0.623 | | 1.1 | 30 |
| Trichloroethene(sim) | 0.483 | 0.434 | | 10.1 | 30 |
| 1,4-Dioxane(sim) | 0.179 | 0.150 | | 16.2 | 30 |
| cis-1,3-Dichloropropene(sim) | 0.498 | 0.453 | | 9.0 | 30 |
| 1,1,2-Trichloroethane(sim) | 0.365 | 0.330 | | 9.6 | 30 |
| Dibromochloromethane(sim) | 0.764 | 0.795 | | -4.1 | 30 |
| 1,2-Dibromoethane(EDB)(sim) | 0.597 | 0.491 | | 17.8 | 30 |
| Tetrachloroethene(sim) | 0.671 | 0.646 | | 3.7 | 30 |
| Bromoform(sim) | 1.657 | 1.852 | | -11.8 | 30 |
| m,p-Xylene(sim) | 2.191 | 2.043 | | 6.8 | 30 |
| 1,1,2,2-Tetrachloroethane(sim) | 1.830 | 1.558 | | 14.9 | 30 |
| Benzyl chloride(sim) | 1.890 | 1.819 | | 3.8 | 30 |
| 1,3-Dichlorobenzene(sim) | 2.412 | 2.175 | | 9.8 | 30 |
| 1,4-Dichlorobenzene(sim) | 2.191 | 1.815 | | 17.2 | 30 |
| sec-Butylbenzene(sim) | 4.557 | 4.106 | | 9.9 | 30 |
| 4-Isopropyltoluene(sim) | 3.859 | 3.549 | | 8.0 | 30 |
| 1,2-Dichlorobenzene(sim) | 2.373 | 2.002 | | 15.6 | 30 |
| n-Butylbenzene(sim) | 3.088 | 2.829 | | 8.4 | 30 |
| 1,2,4-Trichlorobenzene(sim) | 1.595 | 1.047 | | 34.4 # | 30 |
| Hexachlorobutadiene(sim) | 2.397 | 1.962 | | 18.1 | 30 |
| % Bromofluorobenzene | 1.395 | 1.388 | | 0.5 | 30 |
| | | | | | |
| | | | | | |
| | | | | | |

(*) Recommended RRF not met (+) %D exceeds criteria % (#) %D exceeds (maximum) criteria
 %D: 20% of target compounds are allowed to be above criteria %, but must be less than the (maximum) %D
 (#) Maximum %D not met.

7A
AIR CONTINUING CALIBRATION CHECK

Lab Name: Phoenix Environmental Labs Client: _____
 Lab Code: Phoenix Case No.: _____ SAS No.: _____ SDG No.: GCM85136
 Instrument: CHEM39 Calibration Date: 11/14/22 Time: 14:40
 Lab File Id: 1114_02.D Init. Calib. Date(s): 10/30/22 10/31/22
 Heated Purge (Y/N): Y Init. Calib. Times: 18:32 11:56
 GC Column: RTX-1 ; #10157 Method File: 39_AIR_1030.M

| COMPOUND | RRF | RRF1 | RRF MIN | %D | % D LIMITS |
|-------------------------------|-------|-------|---------|---------|------------|
| Propylene | 1.149 | 1.410 | | -22.7 | 30 |
| Dichlorodifluoromethane | 3.820 | 3.861 | | -1.1 | 30 |
| Chloromethane | 2.622 | 3.211 | | -22.5 | 30 |
| 1,2-Dichlorotetrafluoroethane | 4.418 | 4.680 | | -5.9 | 30 |
| Vinyl Chloride | 2.112 | 2.250 | | -6.5 | 30 |
| 1,3-Butadiene | 2.131 | 2.046 | | 4.0 | 30 |
| Bromomethane | 1.638 | 1.590 | | 2.9 | 30 |
| Chloroethane | 0.864 | 0.963 | | -11.5 | 30 |
| Ethanol | 0.886 | 1.088 | | -22.8 | 30 |
| Acetone | 4.178 | 5.660 | | -35.5 # | 30 |
| Trichlorofluoromethane | 4.334 | 4.344 | | -0.2 | 30 |
| Isopropylalcohol | 5.157 | 5.761 | | -11.7 | 30 |
| Acrylonitrile | 2.062 | 2.288 | | -11.0 | 30 |
| 1,1-Dichloroethene | 3.591 | 3.775 | | -5.1 | 30 |
| Methylene Chloride | 3.689 | 4.140 | | -12.2 | 30 |
| Carbon Disulfide | 4.550 | 4.889 | | -7.5 | 30 |
| Trichlorotrifluoroethane | 3.544 | 3.455 | | 2.5 | 30 |
| Trans-1,2-Dichloroethene | 1.832 | 1.960 | | -7.0 | 30 |
| 1,1-Dichloroethane | 2.169 | 2.369 | | -9.2 | 30 |
| Methyl tert-butyl ether(MTBE) | 2.649 | 2.628 | | 0.8 | 30 |
| Methyl Ethyl Ketone | 3.111 | 3.596 | | -15.6 | 30 |
| Cis-1,2-Dichloroethene | 1.781 | 1.918 | | -7.7 | 30 |
| Hexane | 2.248 | 2.363 | | -5.1 | 30 |
| Chloroform | 2.343 | 2.415 | | -3.1 | 30 |
| Ethyl acetate | 0.538 | 0.497 | | 7.6 | 30 |
| Tetrahydrofuran | 1.820 | 2.106 | | -15.7 | 30 |
| 1,2-Dichloroethane | 1.805 | 1.944 | | -7.7 | 30 |
| 1,1,1-Trichloroethane | 2.361 | 2.418 | | -2.4 | 30 |
| Benzene | 3.067 | 3.363 | | -9.7 | 30 |
| Carbon Tetrachloride | 2.540 | 2.435 | | 4.1 | 30 |
| Cyclohexane | 1.218 | 1.262 | | -3.6 | 30 |
| 1,2-dichloropropane | 0.417 | 0.432 | | -3.6 | 30 |
| Bromodichloromethane | 0.722 | 0.642 | | 11.1 | 30 |
| Trichloroethene | 0.474 | 0.419 | | 11.6 | 30 |
| 1,4-Dioxane | 0.181 | 0.141 | | 22.1 | 30 |

(*) Recommended RRF not met (+) %D exceeds criteria % (#) %D exceeds (maximum) criteria
 %D: 20% of target compounds are allowed to be above criteria %, but must be less than the (maximum) %D
 (#) Maximum %D not met.

3
AIR ICV RECOVERY

Lab Name: Phoenix Environmental Labs Client: ACT

Lab Code: Phoenix Case No: _____ SAS No: _____ SDG No GCM85136

ICV - Client Id: ICV_CHEM20_1010

| COMPOUND | SPIKE ADDED (ppbv) | | ICV CONCENTRATION (ppbv) | ICV % REC # | QC. LIMITS REC. |
|-------------------------------|--------------------------|--|--------------------------------|-------------------|-----------------------|
| Propylene | 10 | | 9.268 | 93 | 70 130 |
| Dichlorodifluoromethane | 10 | | 9.484 | 95 | 70 130 |
| Chloromethane | 10 | | 8.908 | 89 | 70 130 |
| 1,2-Dichlorotetrafluoroethane | 10 | | 9.279 | 93 | 70 130 |
| Vinyl Chloride | 10 | | 9.232 | 92 | 70 130 |
| 1,3-Butadiene | 10 | | 9.443 | 94 | 70 130 |
| Bromomethane | 10 | | 9.238 | 92 | 70 130 |
| Chloroethane | 10 | | 9.545 | 95 | 70 130 |
| Ethanol | 10 | | 10.94 | 109 | 70 130 |
| Acetone | 10 | | 9.011 | 90 | 70 130 |
| Trichlorofluoromethane | 10 | | 9.247 | 92 | 70 130 |
| Isopropylalcohol | 10 | | 9.940 | 99 | 70 130 |
| Acrylonitrile | 10 | | 9.946 | 99 | 70 130 |
| 1,1-Dichloroethene | 10 | | 9.536 | 95 | 70 130 |
| Methylene Chloride | 10 | | 8.938 | 89 | 70 130 |
| Carbon Disulfide | 10 | | 9.571 | 96 | 70 130 |
| Trichlorotrifluoroethane | 10 | | 9.599 | 96 | 70 130 |
| Trans-1,2-Dichloroethene | 10 | | 9.363 | 94 | 70 130 |
| 1,1-Dichloroethane | 10 | | 9.472 | 95 | 70 130 |
| Methyl tert-butyl ether(MTBE) | 10 | | 9.505 | 95 | 70 130 |
| Methyl Ethyl Ketone | 10 | | 9.082 | 91 | 70 130 |
| Cis-1,2-Dichloroethene | 10 | | 9.583 | 96 | 70 130 |
| Hexane | 10 | | 9.545 | 95 | 70 130 |
| Chloroform | 10 | | 9.437 | 94 | 70 130 |
| Ethyl acetate | 10 | | 8.240 | 82 | 70 130 |
| Tetrahydrofuran | 10 | | 9.711 | 97 | 70 130 |
| 1,2-Dichloroethane | 10 | | 9.518 | 95 | 70 130 |
| 1,1,1-Trichloroethane | 10 | | 9.031 | 90 | 70 130 |
| Benzene | 10 | | 9.400 | 94 | 70 130 |
| Carbon Tetrachloride | 10 | | 9.484 | 95 | 70 130 |
| Cyclohexane | 10 | | 8.882 | 89 | 70 130 |
| 1,2-dichloropropane | 10 | | 9.554 | 96 | 70 130 |
| Bromodichloromethane | 10 | | 9.565 | 96 | 70 130 |
| Trichloroethene | 10 | | 9.491 | 95 | 70 130 |
| 1,4-Dioxane | 10 | | 9.136 | 91 | 70 130 |
| Heptane | 10 | | 9.591 | 96 | 70 130 |
| cis-1,3-Dichloropropene | 10 | | 9.685 | 97 | 70 130 |
| 4-Methyl-2-pentanone(MIBK) | 10 | | 9.672 | 97 | 70 130 |
| trans-1,3-Dichloropropene | 10 | | 9.522 | 95 | 70 130 |
| 1,1,2-Trichloroethane | 10 | | 9.589 | 96 | 70 130 |
| Toluene | 10 | | 9.536 | 95 | 70 130 |
| Dibromochloromethane | 10 | | 9.646 | 96 | 70 130 |
| 2-Hexanone(MBK) | 10 | | 10.12 | 101 | 70 130 |
| 1,2-Dibromoethane(EDB) | 10 | | 9.427 | 94 | 70 130 |

Data Path : H:\AIR2022\CHEM20\10OCT\10\
 Data File : 1010_17.D
 Acq On : 10 Oct 2022 6:47 pm
 Operator :
 Sample : 10ppbv LCS ; AirT01502A
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Oct 11 09:05:48 2022
 Quant Method : H:\AIR2022\CHEM20\Methods\20_AIR_1010.M
 Quant Title : VOA Standards for 5 point calibration
 QLast Update : Tue Oct 11 09:04:50 2022
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|----------------|----------|--------|--------|----------|
| Internal Standards | | | | | | |
| 1) Bromochloromethane | 7.528 | 130 | 258212 | 10.000 | ng | 0.00 |
| 37) 1,4-Difluorobenzene | 8.682 | 114 | 953852 | 10.000 | ng | 0.00 |
| 54) Chlorobenzene-d5 | 11.135 | 82 | 458458 | 10.000 | ng | 0.00 |
| 81) Bromochloromethane(sim) | 7.533 | 130 | 280275 | 10.000 | ng | # 0.00 |
| 96) 1,4-Difluorobenzene(sim) | 8.682 | 114 | 953423 | 10.000 | ng | 0.00 |
| 106) Chlorobenzene-d5(sim) | 11.135 | 82 | 458458 | 10.000 | ng | 0.00 |
| System Monitoring Compounds | | | | | | |
| 63) % Bromofluorobenzene | 11.966 | 95 | 635979 | 9.942 | ppbv | 0.00 |
| Spiked Amount | 10.000 | Range 70 - 130 | Recovery | = | 99.40% | |
| Target Compounds | | | | | | |
| | | | | | | Qvalue |
| 2) Propylene | 4.078 | 41 | 318948 | 9.268 | ppbv | 99 |
| 3) Dichlorodifluoromethane | 4.153 | 85 | 660568 | 9.484 | ppbv | 99 |
| 4) Chloromethane | 4.304 | 50 | 324553 | 8.907 | ppbv | 100 |
| 5) 1,2-Dichlorotetrafluor... | 4.401 | 85 | 564565 | 9.279 | ppbv | 99 |
| 6) Vinyl Chloride | 4.509 | 62 | 252133 | 9.232 | ppbv | 99 |
| 7) 1,3-Butadiene | 4.638 | 54 | 249476 | 9.443 | ppbv | 97 |
| 8) Bromomethane | 4.865 | 94 | 200710 | 9.238 | ppbv | 100 |
| 9) Chloroethane | 5.005 | 64 | 104631 | 9.545 | ppbv | 98 |
| 10) Vinyl Bromide | 5.285 | 106 | 207923 | 9.537 | ppbv | 99 |
| 11) Ethanol | 5.123 | 45 | 136121 | 10.943 | ppbv | 98 |
| 12) Acetone | 5.501 | 43 | 553064 | 9.011 | ppbv | 100 |
| 13) Trichlorofluoromethane | 5.598 | 101 | 645528 | 9.247 | ppbv | 99 |
| 14) Isopropylalcohol | 5.695 | 45 | 747444 | 9.940 | ppbv | 100 |
| 15) Acrylonitrile | 5.803 | 53 | 276422 | 9.946 | ppbv | 97 |
| 16) 1,1-Dichloroethene | 6.055 | 61 | 439032 | 9.536 | ppbv | 99 |
| 17) Methylene Chloride | 6.124 | 49 | 427087 | 8.938 | ppbv | 99 |
| 18) tert-butyl alcohol | 6.158 | 59 | 658293 | 9.893 | ppbv# | 91 |
| 19) Allyl Chloride | 6.210 | 41 | 472677 | 7.628 | pbv # | 85 |
| 20) Carbon Disulfide | 6.356 | 76 | 629039 | 9.571 | ppbv | 99 |
| 21) Trichlorotrifluoroethane | 6.296 | 101 | 543597 | 9.599 | ppbv | 99 |
| 22) Trans-1,2-Dichloroethene | 6.752 | 61 | 436828 | 9.363 | ppbv | 99 |
| 23) 1,1-Dichloroethane | 6.886 | 63 | 516221 | 9.471 | ppbv | 100 |
| 24) Methyl tert-butyl ethe... | 6.941 | 73 | 622732 | 9.505 | ppbv | 99 |
| 25) Vinyl Acetate | 6.965 | 43 | 2078705 | 10.098 | ppbv# | 100 |
| 26) Methyl Ethyl Ketone | 7.146 | 43 | 827175 | 9.082 | ppbv | 95 |
| 27) Cis-1,2-Dichloroethene | 7.413 | 61 | 419244 | 9.583 | ppbv | 98 |
| 28) Hexane | 7.528 | 57 | 521177 | 9.545 | ppbv | 98 |
| 29) Chloroform | 7.601 | 83 | 525383 | 9.437 | ppbv | 100 |
| 30) Ethyl acetate | 7.538 | 61 | 100682 | 8.240 | ppbv | 99 |
| 31) Tetrahydrofuran | 7.871 | 42 | 468709 | 9.711 | ppbv | 99 |
| 32) 1,2-Dichloroethane | 8.069 | 62 | 406053 | 9.518 | ppbv | 99 |
| 33) 1,1,1-Trichloroethane | 8.226 | 97 | 558446 | 9.031 | ppbv | 99 |
| 34) Benzene | 8.501 | 78 | 712260 | 9.400 | ppbv | 100 |
| 35) Carbon Tetrachloride | 8.592 | 117 | 645644 | 9.484 | ppbv | 99 |
| 36) Cyclohexane | 8.671 | 84 | 272087 | 8.882 | ppbv | 99 |
| 38) 1,2-dichloropropane | 8.977 | 63 | 356284 | 9.554 | ppbv | 98 |
| 39) Bromodichloromethane | 9.090 | 83 | 557079 | 9.565 | ppbv | 99 |
| 40) Trichloroethene | 9.112 | 130 | 401243 | 9.491 | ppbv | 99 |
| 41) 2,2,4-trimethylpentane | 9.112 | 57 | 1797926 | 10.211 | ppbv | 99 |
| 42) 1,4-Dioxane | 9.124 | 88 | 150779 | 9.135 | ppbv | 97 |
| 43) Methyl methacrylate | 9.180 | 69 | 261010 | 9.756 | ppbv | 99 |
| 44) Heptane | 9.237 | 43 | 780278 | 9.591 | ppbv | 100 |

Data Path : H:\AIR2022\CHEM20\10OCT\10\
 Data File : 1010_17.D
 Acq On : 10 Oct 2022 6:47 pm
 Operator :
 Sample : 10ppbv LCS ; AirT01502A
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Oct 11 09:05:48 2022
 Quant Method : H:\AIR2022\CHEM20\Methods\20_AIR_1010.M
 Quant Title : VOA Standards for 5 point calibration
 QLast Update : Tue Oct 11 09:04:50 2022
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------------------|--------|------|----------|--------|-------|----------|
| 45) cis-1,3-Dichloropropene | 9.588 | 75 | 422018 | 9.685 | ppbv | 100 |
| 46) 4-Methyl-2-pentanone(M... | 9.610 | 43 | 1031127 | 9.672 | ppbv | 99 |
| 47) trans-1,3-Dichloropropene | 9.871 | 75 | 395648 | 9.522 | ppbv | 99 |
| 48) 1,1,2-Trichloroethane | 9.984 | 97 | 324351 | 9.589 | ppbv | 99 |
| 49) Toluene | 10.154 | 91 | 931305 | 9.536 | ppbv | 99 |
| 50) Dibromochloromethane | 10.401 | 129 | 685851 | 9.646 | ppbv | 99 |
| 51) 2-Hexanone(MBK) | 10.278 | 43 | 1026154 | 10.123 | ppbv | 98 |
| 52) 1,2-Dibromoethane(EDB) | 10.547 | 107 | 546309 | 9.427 | ppbv | 100 |
| 53) Tetrachloroethene | 10.780 | 166 | 574311 | 10.272 | ppbv | 99 |
| 55) 1,1,1,2-Tetrachloroethane | 11.146 | 131 | 516911 | 10.372 | ppbv | 98 |
| 56) Chlorobenzene | 11.156 | 112 | 792762 | 9.284 | ppbv | 100 |
| 57) Ethylbenzene | 11.351 | 91 | 1202862 | 9.485 | ppbv | 99 |
| 58) m,p-Xylene | 11.443 | 91 | 1878835 | 18.924 | ppbv | 99 |
| 59) Bromoform | 11.525 | 173 | 616070 | 9.080 | ppbv | 99 |
| 60) Styrene | 11.648 | 104 | 739504 | 9.428 | ppbv | 99 |
| 61) 1,1,2,2-Tetrachloroethane | 11.699 | 83 | 664216 | 8.848 | ppbv | 98 |
| 62) o-Xylene | 11.709 | 91 | 969715 | 9.409 | ppbv | 100 |
| 64) 2-Chlorotoluene | 12.314 | 126 | 436987 | 10.385 | ppbv | 97 |
| 65) Isopropylbenzene | 12.027 | 105 | 1552539 | 10.205 | ppbv | 100 |
| 66) n-Propylbenzene | 12.314 | 120 | 460659 | 10.435 | ppbv | 99 |
| 67) 4-Ethyltoluene | 12.396 | 105 | 1430058 | 9.195 | ppbv | 99 |
| 68) 1,3,5-Trimethylbenzene | 12.427 | 105 | 1221443 | 9.598 | ppbv | 100 |
| 69) 1,2,4-Trimethylbenzene | 12.684 | 105 | 1230069 | 9.425 | ppbv | 99 |
| 70) tert-butylbenzene | 12.684 | 119 | 1532392 | 10.249 | ppbv | 99 |
| 71) Benzyl chloride | 12.776 | 91 | 888966 | 9.159 | ppbv | 100 |
| 72) 1,3-Dichlorobenzene | 12.797 | 146 | 846499 | 8.900 | ppbv | 100 |
| 73) 1,4-Dichlorobenzene | 12.838 | 146 | 791252 | 8.816 | ppbv | 99 |
| 74) sec-Butylbenzene | 12.848 | 105 | 1941435 | 10.098 | ppbv | 100 |
| 75) 4-Isopropyltoluene | 12.930 | 119 | 1848420 | 9.888 | ppbv | 100 |
| 76) 1,2-Dichlorobenzene | 13.053 | 146 | 781783 | 8.687 | ppbv | 98 |
| 77) n-Butylbenzene | 13.197 | 91 | 1445038 | 9.752 | ppbv | 99 |
| 78) 1,2,4-Trichlorobenzene | 14.233 | 180 | 575051 | 10.207 | ppbv | 97 |
| 79) Naphthalene | 14.336 | 128 | 1372879 | 10.253 | ppbv# | 100 |
| 80) Hexachlorobutadiene | 14.572 | 225 | 702536 | 10.218 | ppbv | 99 |
| 82] 1,2-Dichlorotetrafluor... | 4.401 | 85 | 564199 | 9.126 | ppbv | 99 |
| 83] Vinyl Chloride(sim) | 4.515 | 62 | 270326 | 8.822 | ppbv | 100 |
| 84] Bromomethane(sim) | 4.865 | 94 | 200710 | 9.205 | ppbv | 100 |
| 85] Trichlorofluoromethane... | 5.603 | 101 | 686376 | 9.959 | ppbv# | 100 |
| 86] 1,2-Dichloroethane(sim) | 8.069 | 62 | 406053 | 9.438 | ppbv | 99 |
| 87] 1,1,1-Trichloroethane(... | 8.231 | 97 | 624529 | 9.908 | ppbv# | 100 |
| 88] Benzene(sim) | 8.501 | 78 | 712877 | 8.915 | ppbv | 100 |
| 89] Carbon Tetrachloride(sim) | 8.597 | 117 | 682498 | 10.417 | ppbv | 100 |
| 90] 1,1-Dichloroethene(sim) | 6.055 | 61 | 439032 | 8.937 | ppbv | 99 |
| 91] Trichlorotrifluoroetha... | 6.302 | 101 | 581716 | 9.929 | ppbv# | 100 |
| 92] Trans-1,2-Dichloroethe... | 6.752 | 61 | 436828 | 9.029 | ppbv | 99 |
| 93] 1,1-Dichloroethane(sim) | 6.891 | 63 | 557272 | 9.611 | ppbv | 100 |
| 94] Cis-1,2-Dichloroethene... | 7.413 | 61 | 419244 | 9.335 | ppbv | 98 |
| 95] Chloroform(sim) | 7.606 | 83 | 554679 | 9.618 | ppbv | 100 |
| 97] 1,2-dichloropropane(sim) | 8.982 | 63 | 382841 | 8.870 | ppbv | 100 |
| 98] Bromodichloromethane(sim) | 9.090 | 83 | 557079 | 9.270 | ppbv | 99 |
| 99] Trichloroethene(sim) | 9.118 | 130 | 435625 | 9.466 | ppbv | 100 |
| 100] 1,4-Dioxane(sim) | 9.124 | 88 | 150779 | 8.820 | ppbv | 97 |
| 101] cis-1,3-Dichloropropen... | 9.593 | 75 | 463845 | 9.775 | ppbv | 100 |
| 102] 1,1,2-Trichloroethane(... | 9.984 | 97 | 324351 | 9.313 | ppbv | 99 |
| 103] Dibromochloromethane(sim) | 10.407 | 129 | 748409 | 10.279 | ppbv | 100 |

Data Path : H:\AIR2022\CHEM20\10OCT\10\
 Data File : 1010_17.D
 Acq On : 10 Oct 2022 6:47 pm
 Operator :
 Sample : 10ppbv LCS ; AirT01502A
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

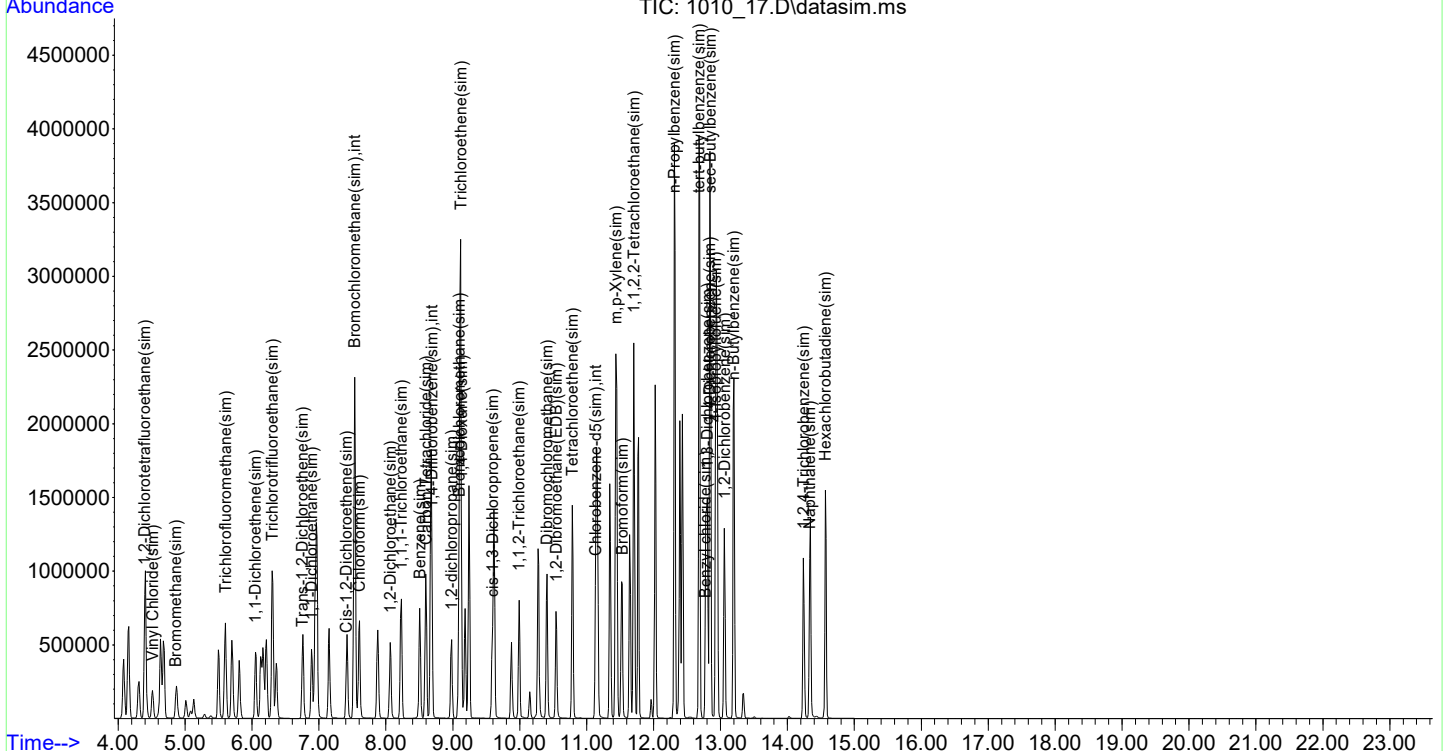
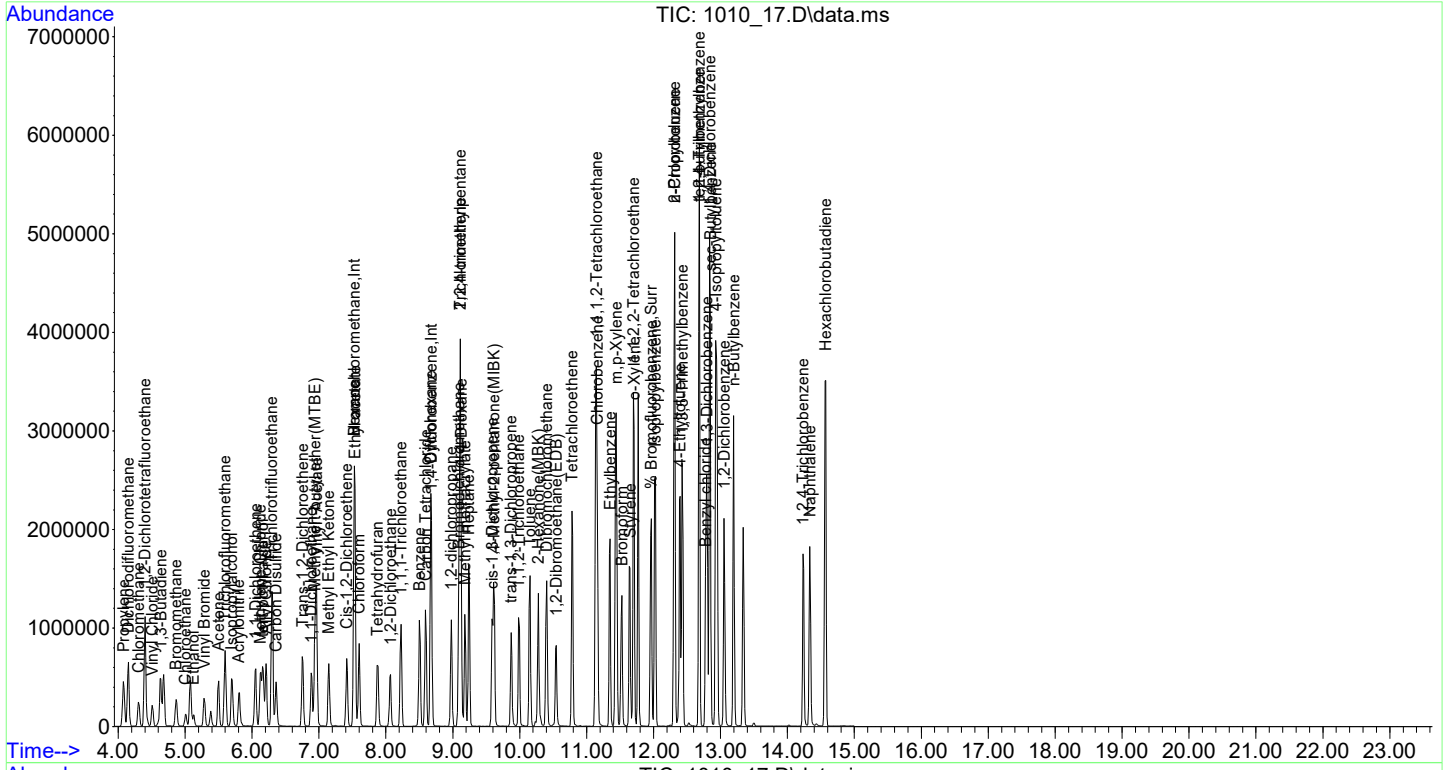
Quant Time: Oct 11 09:05:48 2022
 Quant Method : H:\AIR2022\CHEM20\Methods\20_AIR_1010.M
 Quant Title : VOA Standards for 5 point calibration
 QLast Update : Tue Oct 11 09:04:50 2022
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------------------|--------|------|----------|--------|-------|----------|
| 104] 1,2-Dibromoethane(EDB)... | 10.547 | 107 | 546309 | 9.604 | ppbv | 100 |
| 105] Tetrachloroethene(sim) | 10.785 | 166 | 661008 | 10.331 | ppbv | 100 |
| 107] Bromoform(sim) | 11.531 | 173 | 719071 | 9.465 | ppbv | 100 |
| 108] m,p-Xylene(sim) | 11.443 | 91 | 1896716 | 18.886 | ppbv | 99 |
| 109] 1,1,2,2-Tetrachloroeth... | 11.705 | 83 | 702780 | 8.377 | ppbv | 100 |
| 110] n-Propylbenzene(sim) | 12.314 | 120 | 460659 | 10.359 | ppbv | 99 |
| 111] tert-butylbenzene(sim) | 12.679 | 119 | 1658202 | 10.163 | ppbv | 99 |
| 112] Benzyl chloride(sim) | 12.776 | 91 | 888966 | 10.261 | ppbv | 100 |
| 113] 1,3-Dichlorobenzene(sim) | 12.802 | 146 | 964642 | 8.724 | ppbv | 97 |
| 114] 1,4-Dichlorobenzene(sim) | 12.838 | 146 | 791252 | 7.878 | ppbv | 99 |
| 115] sec-Butylbenzene(sim) | 12.843 | 105 | 2122584 | 10.159 | ppbv | 100 |
| 116] 4-Isopropyltoluene(sim) | 12.930 | 119 | 1848725 | 10.449 | ppbv | 100 |
| 117] 1,2-Dichlorobenzene(sim) | 13.059 | 146 | 869786 | 7.993 | ppbv | 100 |
| 118] n-Butylbenzene(sim) | 13.197 | 91 | 1445038 | 10.209 | ppbv | 99 |
| 119] 1,2,4-Trichlorobenzene... | 14.239 | 180 | 662104 | 9.055 | ppbv | 100 |
| 120] Naphthalene(sim) | 14.336 | 128 | 1372879 | 9.714 | ppbv | 100 |
| 121] Hexachlorobutadiene(sim) | 14.567 | 225 | 876075 | 7.972 | ppbv | 100 |

(#)out of range (m)manual integration reviewed by analyst (+)signals summed

Data Path : H:\AIR2022\CHEM20\10OCT\10\
Data File : 1010_17.D
Acq On : 10 Oct 2022 6:47 pm
Operator :
Sample : 10ppbv LCS ; AirT01502A
Misc :
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Oct 11 09:05:48 2022
Quant Method : H:\AIR2022\CHEM20\Methods\20_AIR_1010.M
Quant Title : VOA Standards for 5 point calibration
QLast Update : Tue Oct 11 09:04:50 2022
Response via : Initial Calibration



ATTACHMENT D

ANNOTATED SUMMARY FORMS

1
AIR ANALYSIS DATA SHEET

CLIENT ID

SS-12

| | | | |
|--------------|-----------------|------------------|--------------------------|
| Client: | <u>ACT</u> | Lab: | <u>Phoenix Env. Labs</u> |
| SDG No.: | <u>GCM85136</u> | Lab Sample ID: | <u>CM85136</u> |
| Canister: | <u>11287</u> | Lab File ID: | <u>1114_34.D</u> |
| Instrument: | <u>CHEM20</u> | Column: | <u>RTX-1 60M</u> |
| | | Date Received: | <u>11/14/22</u> |
| Purge Volume | <u>200</u> (cc) | Date Analyzed: | <u>11/15/22</u> |
| Matrix: | <u>AIR</u> | Dilution Factor: | <u>1</u> |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|-------------|-------------------------------|-------|----|-------|-------|---|
| 115-07-1 | Propylene | 1.08 | J | 0.581 | 0.581 | r |
| 75-71-8 | Dichlorodifluoromethane | 0.481 | J | 0.202 | 0.202 | r |
| 74-87-3 | Chloromethane | 0.485 | UJ | 0.485 | 0.485 | r |
| 106-99-0 | 1,3-Butadiene | 0.452 | UJ | 0.452 | 0.452 | r |
| 75-00-3 | Chloroethane | 0.379 | UJ | 0.379 | 0.379 | r |
| 64-17-5 | Ethanol | 4.22 | J | 0.531 | 0.531 | r |
| 67-64-1 | Acetone | 52.7 | J | 0.421 | 0.421 | r |
| 75-69-4 | Trichlorofluoromethane | 0.237 | J | 0.178 | 0.178 | r |
| 67-63-0 | Isopropylalcohol | 1.07 | J | 0.407 | 0.407 | r |
| 107-13-1 | Acrylonitrile | 0.461 | UJ | 0.461 | 0.461 | r |
| 75-09-2 | Methylene Chloride | 0.863 | UJ | 0.863 | 0.863 | r |
| 75-15-0 | Carbon Disulfide | 2.08 | J | 0.321 | 0.321 | r |
| 1634-04-4 | Methyl tert-butyl ether(MTBE) | 0.278 | UJ | 0.278 | 0.278 | r |
| 78-93-3 | Methyl Ethyl Ketone | 2.13 | J | 0.339 | 0.339 | r |
| 110-54-3 | Hexane | 0.864 | J | 0.284 | 0.284 | r |
| 67-66-3 | Chloroform | 5.41 | J | 0.205 | 0.205 | r |
| 141-78-6 | Ethyl acetate | 0.278 | UJ | 0.278 | 0.278 | r |
| 109-99-9 | Tetrahydrofuran | 0.339 | UJ | 0.339 | 0.339 | r |
| 71-43-2 | Benzene | 0.313 | UJ | 0.313 | 0.313 | r |
| 110-82-7 | Cyclohexane | 0.291 | UJ | 0.291 | 0.291 | r |
| 142-82-5 | Heptane | 0.779 | J | 0.244 | 0.244 | r |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | 0.486 | J | 0.244 | 0.244 | r |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.221 | UJ | 0.221 | 0.221 | r |
| 108-88-3 | Toluene | 0.719 | J | 0.266 | 0.266 | r |
| 591-78-6 | 2-Hexanone(MBK) | 0.244 | UJ | 0.244 | 0.244 | r |
| 127-18-4 | Tetrachloroethene | 5.45 | J | 0.037 | 0.037 | r |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.146 | UJ | 0.146 | 0.146 | r |
| 108-90-7 | Chlorobenzene | 0.217 | UJ | 0.217 | 0.217 | r |
| 100-41-4 | Ethylbenzene | 0.230 | UJ | 0.230 | 0.230 | r |
| 179601-23-1 | m,p-Xylene | 0.848 | J | 0.230 | 0.230 | r |
| 100-42-5 | Styrene | 0.235 | UJ | 0.235 | 0.235 | r |
| 95-47-6 | o-Xylene | 0.346 | J | 0.230 | 0.230 | r |
| 98-82-8 | Isopropylbenzene | 0.204 | UJ | 0.204 | 0.204 | r |
| 622-96-8 | 4-Ethyltoluene | 0.500 | J | 0.204 | 0.204 | r |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.204 | UJ | 0.204 | 0.204 | r |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.929 | J | 0.204 | 0.204 | r |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

1
AIR ANALYSIS DATA SHEET

CLIENT ID

SS-12

| | | | |
|--------------|----------|------------------|-------------------|
| Client: | ACT | Lab: | Phoenix Env. Labs |
| SDG No.: | GCM85136 | Lab Sample ID: | CM85136 |
| Canister: | 11287 | Lab File ID: | 1114_34.D |
| Instrument: | CHEM20 | Column: | RTX-1 60M |
| Purge Volume | 200 (cc) | Date Received: | 11/14/22 |
| Matrix: | AIR | Dilution Factor: | 1 |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|------------|------------------------------------|-------|----|-------|-------|---|
| 76-14-2 | 1,2-Dichlorotetrafluoroethane(sim) | 0.143 | UJ | 0.143 | 0.143 | r |
| 75-01-4 | Vinyl Chloride(sim) | 0.078 | UJ | 0.078 | 0.078 | r |
| 74-83-9 | Bromomethane(sim) | 0.258 | UJ | 0.258 | 0.258 | r |
| 107-06-2 | 1,2-Dichloroethane(sim) | 0.247 | UJ | 0.247 | 0.247 | r |
| 71-55-6 | 1,1,1-Trichloroethane(sim) | 0.183 | UJ | 0.183 | 0.183 | r |
| 56-23-5 | Carbon Tetrachloride(sim) | 0.098 | J | 0.032 | 0.032 | r |
| 75-35-4 | 1,1-Dichloroethene(sim) | 0.051 | UJ | 0.051 | 0.051 | r |
| 76-13-1 | Trichlorotrifluoroethane(sim) | 0.131 | UJ | 0.131 | 0.131 | r |
| 156-60-5 | Trans-1,2-Dichloroethene(sim) | 0.252 | UJ | 0.252 | 0.252 | r |
| 75-34-3 | 1,1-Dichloroethane(sim) | 0.247 | UJ | 0.247 | 0.247 | r |
| 156-59-2 | Cis-1,2-Dichloroethene(sim) | 0.051 | UJ | 0.051 | 0.051 | r |
| 78-87-5 | 1,2-dichloropropane(sim) | 0.217 | UJ | 0.217 | 0.217 | r |
| 75-27-4 | Bromodichloromethane(sim) | 0.149 | UJ | 0.149 | 0.149 | r |
| 79-01-6 | Trichloroethene(sim) | 0.037 | UJ | 0.037 | 0.037 | r |
| 123-91-1 | 1,4-Dioxane(sim) | 0.278 | UJ | 0.278 | 0.278 | r |
| 10061-01-5 | cis-1,3-Dichloropropene(sim) | 0.221 | UJ | 0.221 | 0.221 | r |
| 79-00-5 | 1,1,2-Trichloroethane(sim) | 0.183 | UJ | 0.183 | 0.183 | r |
| 124-48-1 | Dibromochloromethane(sim) | 0.118 | UJ | 0.118 | 0.118 | r |
| 106-93-4 | 1,2-Dibromoethane(EDB)(sim) | 0.130 | UJ | 0.130 | 0.130 | r |
| 75-25-2 | Bromoform(sim) | 0.097 | UJ | 0.097 | 0.097 | r |
| 79-34-5 | 1,1,1,2-Tetrachloroethane(sim) | 0.146 | UJ | 0.146 | 0.146 | r |
| 100-44-7 | Benzyl chloride(sim) | 0.193 | UJ | 0.193 | 0.193 | r |
| 541-73-1 | 1,3-Dichlorobenzene(sim) | 0.166 | UJ | 0.166 | 0.166 | r |
| 106-46-7 | 1,4-Dichlorobenzene(sim) | 0.166 | UJ | 0.166 | 0.166 | r |
| 135-98-8 | sec-Butylbenzene(sim) | 0.182 | UJ | 0.182 | 0.182 | r |
| 99-87-6 | 4-Isopropyltoluene(sim) | 0.182 | UJ | 0.182 | 0.182 | r |
| 95-50-1 | 1,2-Dichlorobenzene(sim) | 0.166 | UJ | 0.166 | 0.166 | r |
| 104-51-8 | n-Butylbenzene(sim) | 0.182 | UJ | 0.182 | 0.182 | r |
| 120-82-1 | 1,2,4-Trichlorobenzene(sim) | 0.135 | UJ | 0.135 | 0.135 | r |
| 87-68-3 | Hexachlorobutadiene(sim) | 0.094 | UJ | 0.094 | 0.094 | r |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

JR 12/5/2025

1
AIR ANALYSIS DATA SHEET

CLIENT ID

IA-11

Client: ACT Lab: Phoenix Env. Labs

SDG No.: GCM85136 Lab Sample ID: CM85137

Canister: 28614 Lab File ID: 1114_13.D

Instrument: CHEM39 Column: FX-1 ; #10157 Date Received: 11/14/22

Purge Volume 200 (cc) Date Analyzed: 11/14/22

Matrix: AIR Dilution Factor: 1

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|------------|------------------------------------|-------|----|-------|-------|---|
| 115-07-1 | Propylene | 0.581 | UJ | 0.581 | 0.581 | r |
| 75-71-8 | Dichlorodifluoromethane | 0.504 | J | 0.202 | 0.202 | r |
| 74-87-3 | Chloromethane | 0.638 | J | 0.485 | 0.485 | r |
| 106-99-0 | 1,3-Butadiene | 0.452 | UJ | 0.452 | 0.452 | r |
| 75-00-3 | Chloroethane | 0.379 | UJ | 0.379 | 0.379 | r |
| 64-17-5 | Ethanol | 70.0 | J | 0.531 | 0.531 | r |
| 67-64-1 | Acetone | 13.9 | J | 0.421 | 0.421 | r |
| 67-63-0 | Isopropylalcohol | 4.17 | J | 0.407 | 0.407 | r |
| 107-13-1 | Acrylonitrile | 0.461 | UJ | 0.461 | 0.461 | r |
| 75-09-2 | Methylene Chloride | 0.863 | UJ | 0.863 | 0.863 | r |
| 75-15-0 | Carbon Disulfide | 0.321 | UJ | 0.321 | 0.321 | r |
| 1634-04-4 | Methyl tert-butyl ether(MTBE) | 0.278 | UJ | 0.278 | 0.278 | r |
| 78-93-3 | Methyl Ethyl Ketone | 0.369 | J | 0.339 | 0.339 | r |
| 110-54-3 | Hexane | 0.284 | UJ | 0.284 | 0.284 | r |
| 141-78-6 | Ethyl acetate | 3.00 | J | 0.278 | 0.278 | r |
| 109-99-9 | Tetrahydrofuran | 0.339 | UJ | 0.339 | 0.339 | r |
| 71-43-2 | Benzene | 0.318 | J | 0.313 | 0.313 | r |
| 110-82-7 | Cyclohexane | 0.291 | UJ | 0.291 | 0.291 | r |
| 142-82-5 | Heptane | 0.244 | UJ | 0.244 | 0.244 | r |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | 0.244 | UJ | 0.244 | 0.244 | r |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.221 | UJ | 0.221 | 0.221 | r |
| 108-88-3 | Toluene | 0.590 | J | 0.266 | 0.266 | r |
| 591-78-6 | 2-Hexanone(MBK) | 0.244 | UJ | 0.244 | 0.244 | r |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.146 | UJ | 0.146 | 0.146 | r |
| 108-90-7 | Chlorobenzene | 0.217 | UJ | 0.217 | 0.217 | r |
| 100-41-4 | Ethylbenzene | 0.230 | UJ | 0.230 | 0.230 | r |
| 100-42-5 | Styrene | 0.235 | UJ | 0.235 | 0.235 | r |
| 95-47-6 | o-Xylene | 0.230 | UJ | 0.230 | 0.230 | r |
| 98-82-8 | Isopropylbenzene | 0.204 | UJ | 0.204 | 0.204 | r |
| 622-96-8 | 4-Ethyltoluene | 0.204 | UJ | 0.204 | 0.204 | r |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.204 | UJ | 0.204 | 0.204 | r |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.204 | UJ | 0.204 | 0.204 | r |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane(sim) | 0.143 | UJ | 0.143 | 0.143 | r |
| 75-01-4 | Vinyl Chloride(sim) | 0.078 | UJ | 0.078 | 0.078 | r |
| 74-83-9 | Bromomethane(sim) | 0.258 | UJ | 0.258 | 0.258 | r |
| 75-69-4 | Trichlorofluoromethane(sim) | 0.192 | J | 0.178 | 0.178 | r |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

1
AIR ANALYSIS DATA SHEET

CLIENT ID

OA-2

| | | | |
|----------------|----------|------------------|-------------------|
| Client: | ACT | Lab: | Phoenix Env. Labs |
| SDG No.: | GCM85136 | Lab Sample ID: | CM85138 |
| Canister: | 11257 | Lab File ID: | 1114_15.D |
| Instrument: | CHEM39 | Column: | TX-1 ; #10157 |
| Date Received: | 11/14/22 | | |
| Purge Volume | 200 | (cc) | 11/14/22 |
| Date Analyzed: | 11/14/22 | | |
| Matrix: | AIR | Dilution Factor: | 1 |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|------------|------------------------------------|-------|----|-------|-------|---|
| 115-07-1 | Propylene | 0.581 | UJ | 0.581 | 0.581 | r |
| 75-71-8 | Dichlorodifluoromethane | 0.479 | J | 0.202 | 0.202 | r |
| 74-87-3 | Chloromethane | 0.576 | J | 0.485 | 0.485 | r |
| 106-99-0 | 1,3-Butadiene | 0.452 | UJ | 0.452 | 0.452 | r |
| 75-00-3 | Chloroethane | 0.379 | UJ | 0.379 | 0.379 | r |
| 64-17-5 | Ethanol | 14.1 | J | 0.531 | 0.531 | r |
| 67-64-1 | Acetone | 4.29 | J | 0.421 | 0.421 | r |
| 67-63-0 | Isopropylalcohol | 2.29 | J | 0.407 | 0.407 | r |
| 107-13-1 | Acrylonitrile | 0.461 | UJ | 0.461 | 0.461 | r |
| 75-09-2 | Methylene Chloride | 0.863 | UJ | 0.863 | 0.863 | r |
| 75-15-0 | Carbon Disulfide | 0.321 | UJ | 0.321 | 0.321 | r |
| 1634-04-4 | Methyl tert-butyl ether(MTBE) | 0.278 | UJ | 0.278 | 0.278 | r |
| 78-93-3 | Methyl Ethyl Ketone | 0.339 | UJ | 0.339 | 0.339 | r |
| 110-54-3 | Hexane | 0.284 | UJ | 0.284 | 0.284 | r |
| 141-78-6 | Ethyl acetate | 0.278 | UJ | 0.278 | 0.278 | r |
| 109-99-9 | Tetrahydrofuran | 0.339 | UJ | 0.339 | 0.339 | r |
| 71-43-2 | Benzene | 0.313 | UJ | 0.313 | 0.313 | r |
| 110-82-7 | Cyclohexane | 0.291 | UJ | 0.291 | 0.291 | r |
| 142-82-5 | Heptane | 0.244 | UJ | 0.244 | 0.244 | r |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | 0.244 | UJ | 0.244 | 0.244 | r |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.221 | UJ | 0.221 | 0.221 | r |
| 108-88-3 | Toluene | 0.520 | J | 0.266 | 0.266 | r |
| 591-78-6 | 2-Hexanone(MBK) | 0.244 | UJ | 0.244 | 0.244 | r |
| 127-18-4 | Tetrachloroethene | 0.280 | J | 0.037 | 0.037 | r |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.146 | UJ | 0.146 | 0.146 | r |
| 108-90-7 | Chlorobenzene | 0.217 | UJ | 0.217 | 0.217 | r |
| 100-41-4 | Ethylbenzene | 0.230 | UJ | 0.230 | 0.230 | r |
| 100-42-5 | Styrene | 0.235 | UJ | 0.235 | 0.235 | r |
| 95-47-6 | o-Xylene | 0.230 | UJ | 0.230 | 0.230 | r |
| 98-82-8 | Isopropylbenzene | 0.204 | UJ | 0.204 | 0.204 | r |
| 622-96-8 | 4-Ethyltoluene | 0.204 | UJ | 0.204 | 0.204 | r |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.204 | UJ | 0.204 | 0.204 | r |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.204 | UJ | 0.204 | 0.204 | r |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane(sim) | 0.143 | UJ | 0.143 | 0.143 | r |
| 75-01-4 | Vinyl Chloride(sim) | 0.078 | UJ | 0.078 | 0.078 | r |
| 74-83-9 | Bromomethane(sim) | 0.258 | UJ | 0.258 | 0.258 | r |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

1
AIR ANALYSIS DATA SHEET

CLIENT ID

OA-2

| | | | |
|--------------|----------|------------------|-------------------|
| Client: | ACT | Lab: | Phoenix Env. Labs |
| SDG No.: | GCM85136 | Lab Sample ID: | CM85138 |
| Canister: | 11257 | Lab File ID: | 1114_15.D |
| Instrument: | CHEM39 | Column: | TX-1 ; #10157 |
| Purge Volume | 200 (cc) | Date Received: | 11/14/22 |
| Matrix: | AIR | Date Analyzed: | 11/14/22 |
| | | Dilution Factor: | 1 |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|-------------|--------------------------------|-------|----|-------|-------|---|
| 75-69-4 | Trichlorofluoromethane(sim) | 0.195 | J | 0.178 | 0.178 | r |
| 107-06-2 | 1,2-Dichloroethane(sim) | 0.247 | UJ | 0.247 | 0.247 | r |
| 71-55-6 | 1,1,1-Trichloroethane(sim) | 0.183 | UJ | 0.183 | 0.183 | r |
| 56-23-5 | Carbon Tetrachloride(sim) | 0.067 | J | 0.032 | 0.032 | r |
| 75-35-4 | 1,1-Dichloroethene(sim) | 0.051 | UJ | 0.051 | 0.051 | r |
| 76-13-1 | Trichlorotrifluoroethane(sim) | 0.131 | UJ | 0.131 | 0.131 | r |
| 156-60-5 | Trans-1,2-Dichloroethene(sim) | 0.252 | UJ | 0.252 | 0.252 | r |
| 75-34-3 | 1,1-Dichloroethane(sim) | 0.247 | UJ | 0.247 | 0.247 | r |
| 156-59-2 | Cis-1,2-Dichloroethene(sim) | 0.051 | UJ | 0.051 | 0.051 | r |
| 67-66-3 | Chloroform(sim) | 0.205 | UJ | 0.205 | 0.205 | r |
| 78-87-5 | 1,2-dichloropropane(sim) | 0.217 | UJ | 0.217 | 0.217 | r |
| 75-27-4 | Bromodichloromethane(sim) | 0.149 | UJ | 0.149 | 0.149 | r |
| 79-01-6 | Trichloroethene(sim) | 0.055 | J | 0.037 | 0.037 | r |
| 123-91-1 | 1,4-Dioxane(sim) | 0.278 | UJ | 0.278 | 0.278 | r |
| 10061-01-5 | cis-1,3-Dichloropropene(sim) | 0.221 | UJ | 0.221 | 0.221 | r |
| 79-00-5 | 1,1,2-Trichloroethane(sim) | 0.183 | UJ | 0.183 | 0.183 | r |
| 124-48-1 | Dibromochloromethane(sim) | 0.118 | UJ | 0.118 | 0.118 | r |
| 106-93-4 | 1,2-Dibromoethane(EDB)(sim) | 0.130 | UJ | 0.130 | 0.130 | r |
| 75-25-2 | Bromoform(sim) | 0.097 | UJ | 0.097 | 0.097 | r |
| 179601-23-1 | m,p-Xylene(sim) | 0.268 | J | 0.230 | 0.230 | r |
| 79-34-5 | 1,1,2,2-Tetrachloroethane(sim) | 0.146 | UJ | 0.146 | 0.146 | r |
| 100-44-7 | Benzyl chloride(sim) | 0.193 | UJ | 0.193 | 0.193 | r |
| 541-73-1 | 1,3-Dichlorobenzene(sim) | 0.166 | UJ | 0.166 | 0.166 | r |
| 106-46-7 | 1,4-Dichlorobenzene(sim) | 0.166 | UJ | 0.166 | 0.166 | r |
| 135-98-8 | sec-Butylbenzene(sim) | 0.182 | UJ | 0.182 | 0.182 | r |
| 99-87-6 | 4-Isopropyltoluene(sim) | 0.182 | UJ | 0.182 | 0.182 | r |
| 95-50-1 | 1,2-Dichlorobenzene(sim) | 0.166 | UJ | 0.166 | 0.166 | r |
| 104-51-8 | n-Butylbenzene(sim) | 0.182 | UJ | 0.182 | 0.182 | r |
| 120-82-1 | 1,2,4-Trichlorobenzene(sim) | 0.135 | UJ | 0.135 | 0.135 | r |
| 87-68-3 | Hexachlorobutadiene(sim) | 0.094 | UJ | 0.094 | 0.094 | r |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |

FORM I AIR
r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

JR 12/5/2025

December 8, 2025

Mr. Jason Stewart
Advanced Cleanup Technologies
228 Park Ave S PMB 34864
New York, New York 10003

Re: Data Usability Summary Report – Phoenix Environmental Laboratories, Inc. – GCO73252 V1

Dear Mr. Stewart:

The evaluation of analytical data by Phoenix Environmental Laboratories for project 9268-BXNY, which were reported in a single data package under Job No. GCO73252 V1 has been completed. The following samples were reported.

IA-4 IA-3 IA-7

Analysis was performed in accordance with EPA Method TO-15 (volatile organics). The review was performed to the extent possible, in accordance with the analytical method, and “DER-10/ Technical Guidance for Site Investigation and Remediation”. Professional judgment is applied as necessary and appropriate. National Functional Guidelines for Organic Data Review was consulted as needed. Qualifiers consistent with those defined by EPA Region 2 are applied as necessary and appropriate.

Below is the Data Usability Summary Report (DUSR) associated with these samples.

| Data Usability Summary Report | |
|---|--------------------------------|
| 1. Is the data package complete as defined under the requirements for the most current DEC ASP Category B or USEPA CLP data deliverables? | No – see Documentation Section |
| 2. Have all holding times been met? | Yes |
| 3. Do all the QC data; blanks, instrument tunings, calibration standards, calibration verifications, surrogate recoveries, spike recoveries, replicate analyses, laboratory controls and sample data fall within the protocol required limits and specifications? | No -see following sections |
| 4. Have all of the data been generated using established and agreed upon analytical protocols? | Yes |
| 5. Does an evaluation of the raw data confirm the results provided in the data summary sheet and the quality control verification forms? | Yes |
| 6. Have the correct data qualifiers been used and are they consistent with the most current DEC ASP? | Yes |

| | |
|---|-----|
| 7. Have any quality control (QC) exceedances been specifically noted in the DUSR and have the corresponding QC summary sheet from the data package been attached to the DUSR? | Yes |
|---|-----|

Overall Evaluation

Based on the data review effort, results are usable, with the following qualifications. For samples that are qualified as estimated (J-, UJ), detected results may be biased low. False negatives may exist in non-detect results. Sample results that are qualified as estimated (J+) may be biased high. For samples that are qualified as estimated with any combination of (J), (J-) and/or (J+), the (J) qualifier takes precedence and is applied to the sample result. It is not possible to determine the direction of the bias and the overall effect on the result.

- The result for ethanol in IA-4 is qualified as estimated (J-) due to low recovery in the laboratory control sample (LCS).
- The results for tetrahydrofuran (undiluted analysis) in IA-3, acetone and tetrahydrofuran (undiluted analysis) in IA-7 are qualified as estimated (J) because the concentration is outside the calibration range.

Qualifier definitions are provided in Attachment A. A copy of the chain of custody record is provided in Attachment B. Pages from the data package illustrating the exceedances and issues described in this validation report are included in Attachment C. Annotated Summary Forms are included in Attachment D detailing qualifications resulting from the data review effort.

The following components were reviewed, where applicable:

- Chain of Custody
- Receiving conditions
- Holding times
- Preservation
- Analyte lists
- Reporting limits
- Requested methods
- Units, and
- Sample related quality control data:
 - Method, instrument blanks
 - Clean canister certification
 - Field blanks
 - Surrogate recoveries
 - LCS recoveries
 - Internal standard area response
 - Duplicates
 - Analyte Identification

- Instrument related quality control data:
 - Instrument tunes
 - Calibration summaries

The following sections of the report detail only quality control exceedances that impacted results. Where a quality control item exceeded control limits but there is no impact to the samples results, these are not detailed in the report.

Documentation: A completeness review of the data package was performed, and the data package was determined to be a complete Category B data package, with the following exceptions:

- The data package does not include summary forms and raw data for the second source initial calibration verification (ICV) standard. The laboratory was contacted and provided the ICV data and is included in attachment C.
- Calibration data for the clean canister certification analysis is not included in the data package.

The following documentation issues were observed during the review:

- Results for ethanol, acetone, isopropyl alcohol, and hexane in all samples are flagged 'S' which the data deliverable identifies this flag as 'This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level.' Where the sample detections are greater than five times the reporting limit, the validator removed the 'S' qualifier.
- The data package documents the sample collection date as August 10, 2023; however, the COC indicates a collection date of August 9, 2023.

Holding Times, Preservation, Sample Integrity:

A copy of the applicable chain of custody (COC) record was included in the data package, documenting sample collection date of August 9, 2023. The samples were received on August 11, 2023. Samples were received intact and analyzed within the method specified hold time.

A. Volatile Organics

1. Laboratory Control Sample (LCS)

Two LCSs were prepared and analyzed with the sample. The LCS is evaluated using control limit of 70-130%R. All recoveries are acceptable, with the following exceptions:

| Analyte | LCS %R | Affected Sample | Qualifier Applied |
|--------------------|-----------|-----------------|----------------------|
| <i>CO71323 LCS</i> | | | |
| Ethanol | 66 | IA-4 | J- |

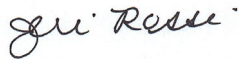
*The summary form provided by the laboratory reports LCS recoveries based on the full scan analysis. The validator reviewed the raw data to determine the recoveries in the SIM analysis and recoveries are acceptable. The result for ethanol in IA-4 is qualified as estimated (J-) due to low recovery in the LCS.

2. Compound Quantitation

The results for tetrahydrofuran (undiluted analysis) in IA-3, acetone and tetrahydrofuran (undiluted analysis) in IA-7 are qualified as estimated (J) because the concentration is outside the calibration range. Results from the diluted analysis should be used.

No other sample results are qualified. Please feel free to contact me at (908) 370-3431 or richjerirossi513@gmail.com if you have any questions regarding this data package review report or need further information.

Sincerely,



Jeri L Rossi, CEAC

Environmental Consulting Chemist

ATTACHMENT A

Qualifier Definitions

EPA Qualifier Definitions

- U The analyte was analyzed for but was not detected above the level of the reported sample quantitation limit.
- J The result is an estimated quantity. The associated numerical value is the approximate concentration of the analyte in the sample.
- J+ The result is an estimated quantity, but the result may be biased high.
- J- The result is an estimated quantity, but the result may be biased low.
- NJ The analyte has been “tentatively identified” or “presumptively” as present and the associated numerical value is the estimated concentration in the sample.
- UJ The analyte was analyzed for but was not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
- R The data are unusable. The sample results are rejected due to serious deficiencies in meeting QC criteria. The analyte may or may not be present in the sample.

ATTACHMENT B

CHAIN OF CUSTODY (COC)



157 East Middle Turnpike, P.O. Box 570, Manchester, CT 06040
 Telephone: 860.645.1102 • Fax: 860.645.0823

ACT

CHAIN OF CUSTODY RECORD

AIR ANALYSES

860-645-1102

email: greg@phoenixlabs.com

P.O. #

Page 1 of 1

Data Delivery:

Fax #: _____
 Email: _____
 Phone #: _____

| | | |
|---|----------------------------|---|
| Report to: Ray Buzeta | Project Name: 9628-BXNY | Data Format: (Circle) Equis Excel Other: |
| Customer: Advanced Cleanup Technologies, Inc. | Invoice to: Karen Friedman | Requested Deliverable: RCP ASP CAT B |
| Address: 228 Park Ave S PMB 34864 | Sampled by: Tim Young | MCP NJ Deliverables |
| 14905 New York, New York 10003 | Quote Number: | |

| Phoenix ID # | Client Sample ID | Canister ID # | Canister Size (L) | Outgoing Canister Pressure ("Hg) | Incoming Canister Pressure ("Hg) | Flow Regulator ID # | Flow Controller Setting (mL/min) | Sampling Start Time | Sampling End Time | Sample Start Date | Canister Pressure at Start ("Hg) | Canister Pressure at End ("Hg) | Ambient/Indoor Air | Soil Gas | Grab (C) Composite (C) | TO-15 | APH |
|-------------------------------|------------------|---------------|-------------------|----------------------------------|----------------------------------|---------------------|----------------------------------|---------------------|-------------------|-------------------|----------------------------------|--------------------------------|--------------------|----------|------------------------|-------|-----|
| THIS SECTION FOR LAB USE ONLY | | | | | | | | | | | | | MATRIX | | ANALYSES | | |
| 73252 | IA-4 | 28570 | 6.0L | -30 | -11 | 2854 | 3.01 | 8/25 | 0945 | 8/23 | -30 | -10 | X | | | X | |
| 73253 | IA-3 | 23348 | 6.0L | -30 | -9 | 2860 | 3.17 | 8/30 | 0910 | 8/23 | -30 | -9 | X | | | X | |
| 73254 | IA-7 | 489 | 6.0L | -30 | -10 | 2988 | 3.28 | 8/20 | 0900 | 8/23 | -30 | -7 | X | | | X | |
| | | | | | | | | ↑ | ↑ | | | | | | | | |
| | | | | | | | | 8/23 | 8/10/23 | | | | | | | | |

| | | | | |
|------------------|--------------|---------|-------|---|
| Relinquished by: | Accepted by: | Date: | Time: | I attest that all media released by Phoenix Environmental Laboratories, Inc. have been received in good working condition and agree to the terms and conditions as listed on the back of this document: |
| | Zu | 8-11-23 | 941 | |
| | Zu | 8-11-23 | 1521 | Signature: [Signature] Date: 8/10/23 |

| | | |
|--|--|--|
| State Where Samples Collected: Bronx, NY | Turnaround Time: <input type="checkbox"/> 1 Day* <input type="checkbox"/> 2 Day* <input type="checkbox"/> 3 Day* <input type="checkbox"/> 4 Day* <input type="checkbox"/> 5 Day* <input checked="" type="checkbox"/> Standard *SURCHARGES MAY APPLY | Requested Criteria: (Please Circle) CT: TAC I/C TAC RES SVVC I/C SVVC RES GWV I/C GWV CES MA: Indoor Air Residential Indoor Air Residential Ind/Commercial Soil Gas Residential Ind/Commercial NJ: Indoor Air Residential Indoor Air Residential Ind/Commercial Soil Gas Residential Ind/Commercial NY: Vapor Intrusion PA: Indoor Air Residential Non-residential VT: Indoor Air Residential Industrial Sub-slab Residential Industrial |
| SPECIAL INSTRUCTIONS, QC REQUIREMENTS, REGULATORY INFORMATION: (3) - 6.0L 24 hr | | |

08/29/2023

Phoenix Environmental Laboratories, Inc.

Page 25 of 297

ATTACHMENT C

**SELECTED PAGES FROM DATA PACKAGE –
QC EXCEEDANCES AND VALIDATION ISSUES**

3
AIR LCS RECOVERY

Lab Name: Phoenix Environmental Labs Client: ACT

Lab Code: Phoenix Case No: _____ SAS No: _____ SDG No GCO73252

LCS - Client Id: CO71323 LCS

| COMPOUND | SPIKE ADDED (ppbv) | | LCS CONCENTRATION (ppbv) | LCS % REC # | QC. LIMITS REC. | |
|-------------------------------|--------------------|--|--------------------------|-------------|-----------------|-----|
| Propylene | 10 | | 9.686 | 97 | 70 | 130 |
| Dichlorodifluoromethane | 10 | | 10.49 | 105 | 70 | 130 |
| Chloromethane | 10 | | 10.12 | 101 | 70 | 130 |
| 1,2-Dichlorotetrafluoroethane | 10 | | 10.44 | 104 | 70 | 130 |
| Vinyl Chloride | 10 | | 10.02 | 100 | 70 | 130 |
| 1,3-Butadiene | 10 | | 10.33 | 103 | 70 | 130 |
| Bromomethane | 10 | | 9.861 | 99 | 70 | 130 |
| Chloroethane | 10 | | 9.859 | 99 | 70 | 130 |
| Ethanol | 10 | | 6.621 | 66 * | 70 | 130 |
| Acetone | 10 | | 9.984 | 100 | 70 | 130 |
| Trichlorofluoromethane | 10 | | 10.02 | 100 | 70 | 130 |
| Isopropylalcohol | 10 | | 9.632 | 96 | 70 | 130 |
| Acrylonitrile | 10 | | 9.573 | 96 | 70 | 130 |
| 1,1-Dichloroethene | 10 | | 9.953 | 100 | 70 | 130 |
| Methylene Chloride | 10 | | 9.926 | 99 | 70 | 130 |
| Carbon Disulfide | 10 | | 9.808 | 98 | 70 | 130 |
| Trichlorotrifluoroethane | 10 | | 10.04 | 100 | 70 | 130 |
| Trans-1,2-Dichloroethene | 10 | | 9.653 | 97 | 70 | 130 |
| 1,1-Dichloroethane | 10 | | 9.760 | 98 | 70 | 130 |
| Methyl tert-butyl ether(MTBE) | 10 | | 10.21 | 102 | 70 | 130 |
| Methyl Ethyl Ketone | 10 | | 9.660 | 97 | 70 | 130 |
| Cis-1,2-Dichloroethene | 10 | | 9.977 | 100 | 70 | 130 |
| Hexane | 10 | | 10.11 | 101 | 70 | 130 |
| Chloroform | 10 | | 10.17 | 102 | 70 | 130 |
| Ethyl acetate | 10 | | 7.794 | 78 | 70 | 130 |
| Tetrahydrofuran | 10 | | 10.70 | 107 | 70 | 130 |
| 1,2-Dichloroethane | 10 | | 10.10 | 101 | 70 | 130 |
| 1,1,1-Trichloroethane | 10 | | 10.10 | 101 | 70 | 130 |
| Benzene | 10 | | 9.960 | 100 | 70 | 130 |
| Carbon Tetrachloride | 10 | | 10.28 | 103 | 70 | 130 |
| Cyclohexane | 10 | | 9.410 | 94 | 70 | 130 |
| 1,2-dichloropropane | 10 | | 9.702 | 97 | 70 | 130 |
| Bromodichloromethane | 10 | | 9.806 | 98 | 70 | 130 |
| Trichloroethene | 10 | | 10.15 | 102 | 70 | 130 |
| 1,4-Dioxane | 10 | | 10.33 | 103 | 70 | 130 |
| Heptane | 10 | | 9.884 | 99 | 70 | 130 |
| cis-1,3-Dichloropropene | 10 | | 10.14 | 101 | 70 | 130 |
| 4-Methyl-2-pentanone(MIBK) | 10 | | 10.13 | 101 | 70 | 130 |
| trans-1,3-Dichloropropene | 10 | | 10.43 | 104 | 70 | 130 |
| 1,1,2-Trichloroethane | 10 | | 9.932 | 99 | 70 | 130 |
| Toluene | 10 | | 10.17 | 102 | 70 | 130 |
| Dibromochloromethane | 10 | | 10.27 | 103 | 70 | 130 |
| 2-Hexanone(MBK) | 10 | | 10.19 | 102 | 70 | 130 |
| 1,2-Dibromoethane(EDB) | 10 | | 10.02 | 100 | 70 | 130 |

FORM III AIR

3
AIR ICV RECOVERY

Lab Name: Phoenix Environmental Labs Client: ACT

Lab Code: Phoenix Case No: _____ SAS No: _____ SDG No GCO73252

ICV - Client Id: ICV_CHEM39_0523

| COMPOUND | SPIKE ADDED (ppbv) | | ICV CONCENTRATION (ppbv) | ICV % REC # | QC. LIMITS REC. | |
|-------------------------------|--------------------|--|--------------------------|-------------|-----------------|-----|
| Propylene | 10 | | 10.46 | 105 | 70 | 130 |
| Dichlorodifluoromethane | 10 | | 10.66 | 107 | 70 | 130 |
| Chloromethane | 10 | | 10.57 | 106 | 70 | 130 |
| 1,2-Dichlorotetrafluoroethane | 10 | | 10.70 | 107 | 70 | 130 |
| Vinyl Chloride | 10 | | 10.63 | 106 | 70 | 130 |
| 1,3-Butadiene | 10 | | 10.77 | 108 | 70 | 130 |
| Bromomethane | 10 | | 10.40 | 104 | 70 | 130 |
| Chloroethane | 10 | | 10.41 | 104 | 70 | 130 |
| Ethanol | 10 | | 9.480 | 95 | 70 | 130 |
| Acetone | 10 | | 10.71 | 107 | 70 | 130 |
| Trichlorofluoromethane | 10 | | 10.31 | 103 | 70 | 130 |
| Isopropylalcohol | 10 | | 8.899 | 89 | 70 | 130 |
| Acrylonitrile | 10 | | 10.20 | 102 | 70 | 130 |
| 1,1-Dichloroethene | 10 | | 10.37 | 104 | 70 | 130 |
| Methylene Chloride | 10 | | 10.39 | 104 | 70 | 130 |
| Carbon Disulfide | 10 | | 10.21 | 102 | 70 | 130 |
| Trichlorotrifluoroethane | 10 | | 10.36 | 104 | 70 | 130 |
| Trans-1,2-Dichloroethene | 10 | | 10.54 | 105 | 70 | 130 |
| 1,1-Dichloroethane | 10 | | 10.39 | 104 | 70 | 130 |
| Methyl tert-butyl ether(MTBE) | 10 | | 10.82 | 108 | 70 | 130 |
| Methyl Ethyl Ketone | 10 | | 10.37 | 104 | 70 | 130 |
| Cis-1,2-Dichloroethene | 10 | | 10.90 | 109 | 70 | 130 |
| Hexane | 10 | | 10.91 | 109 | 70 | 130 |
| Chloroform | 10 | | 10.57 | 106 | 70 | 130 |
| Ethyl acetate | 10 | | 8.404 | 84 | 70 | 130 |
| Tetrahydrofuran | 10 | | 11.91 | 119 | 70 | 130 |
| 1,2-Dichloroethane | 10 | | 10.58 | 106 | 70 | 130 |
| 1,1,1-Trichloroethane | 10 | | 10.41 | 104 | 70 | 130 |
| Benzene | 10 | | 10.61 | 106 | 70 | 130 |
| Carbon Tetrachloride | 10 | | 10.50 | 105 | 70 | 130 |
| Cyclohexane | 10 | | 9.886 | 99 | 70 | 130 |
| 1,2-dichloropropane | 10 | | 10.61 | 106 | 70 | 130 |
| Bromodichloromethane | 10 | | 10.39 | 104 | 70 | 130 |
| Trichloroethene | 10 | | 10.48 | 105 | 70 | 130 |
| 1,4-Dioxane | 10 | | 10.71 | 107 | 70 | 130 |
| Heptane | 10 | | 10.79 | 108 | 70 | 130 |
| cis-1,3-Dichloropropene | 10 | | 10.92 | 109 | 70 | 130 |
| 4-Methyl-2-pentanone(MIBK) | 10 | | 11.01 | 110 | 70 | 130 |
| trans-1,3-Dichloropropene | 10 | | 11.33 | 113 | 70 | 130 |
| 1,1,2-Trichloroethane | 10 | | 10.53 | 105 | 70 | 130 |
| Toluene | 10 | | 10.90 | 109 | 70 | 130 |
| Dibromochloromethane | 10 | | 10.64 | 106 | 70 | 130 |
| 2-Hexanone(MBK) | 10 | | 11.30 | 113 | 70 | 130 |
| 1,2-Dibromoethane(EDB) | 10 | | 10.71 | 107 | 70 | 130 |

Data Path : H:\AIR2023\CHEM39\05MAY\23\
 Data File : 0523_20.D
 Acq On : 24 May 2023 1:15 am
 Operator :
 Sample : 10ppb lcs ; to1503t
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: May 25 16:52:43 2023
 Quant Method : H:\AIR2023\CHEM39\Methods\39_AIR_0523.M
 Quant Title : VOA Standards for 5 point calibration
 QLast Update : Thu May 25 16:51:55 2023
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) | |
|-------------------------------|--------|----------------|----------|--------|---------|----------|--------|
| Internal Standards | | | | | | | |
| 1) Bromochloromethane | 7.521 | 130 | 423878 | 10.000 | ng | 0.00 | |
| 37) 1,4-Difluorobenzene | 8.687 | 114 | 1517717 | 10.000 | ng | 0.00 | |
| 54) Chlorobenzene-d5 | 11.155 | 82 | 720064 | 10.000 | ng | 0.00 | |
| 81) Bromochloromethane(sim) | 7.524 | 130 | 412681 | 10.000 | ng | # 0.00 | |
| 96) 1,4-Difluorobenzene(sim) | 8.687 | 114 | 1517717 | 10.000 | ng | 0.00 | |
| 106) Chlorobenzene-d5(sim) | 11.155 | 82 | 720842 | 10.000 | ng | 0.00 | |
| System Monitoring Compounds | | | | | | | |
| 63) % Bromofluorobenzene | 11.989 | 95 | 1038163 | 10.147 | ppbv | 0.00 | |
| Spiked Amount | 10.000 | Range 70 - 130 | Recovery | = | 101.50% | | |
| Target Compounds | | | | | | | |
| | | | | | | | Qvalue |
| 2) Propylene | 4.033 | 41 | 221135 | 10.463 | ppbv | | 98 |
| 3) Dichlorodifluoromethane | 4.098 | 85 | 1086873 | 10.661 | ppbv | | 99 |
| 4) Chloromethane | 4.252 | 50 | 267694 | 10.574 | ppbv | | 98 |
| 5) 1,2-Dichlorotetrafluor... | 4.349 | 85 | 1110004 | 10.702 | ppbv | | 98 |
| 6) Vinyl Chloride | 4.463 | 62 | 363834 | 10.628 | ppbv | | 100 |
| 7) 1,3-Butadiene | 4.584 | 54 | 245076 | 10.774 | ppbv | | 98 |
| 8) Bromomethane | 4.828 | 94 | 423424 | 10.402 | ppbv | | 99 |
| 9) Chloroethane | 4.965 | 64 | 164381 | 10.411 | ppbv | | 99 |
| 10) Vinyl Bromide | 5.253 | 106 | 416390 | 10.158 | ppbv | | 99 |
| 11) Ethanol | 5.055 | 45 | 75927 | 9.480 | ppbv | | 98 |
| 12) Acetone | 5.437 | 43 | 455073 | 10.710 | ppbv | | 97 |
| 13) Trichlorofluoromethane | 5.565 | 101 | 1120203 | 10.305 | ppbv | | 99 |
| 14) Isopropylalcohol | 5.621 | 45 | 424785 | 8.899 | ppbv | | 97 |
| 15) Acrylonitrile | 5.768 | 53 | 220536 | 10.202 | ppbv | | 99 |
| 16) 1,1-Dichloroethene | 6.025 | 61 | 564995 | 10.373 | ppbv | | 94 |
| 17) Methylene Chloride | 6.103 | 49 | 336078 | 10.385 | ppbv | | 89 |
| 18) tert-butyl alcohol | 6.067 | 59 | 613038 | 10.187 | ppbv | | 99 |
| 19) Allyl Chloride | 6.186 | 41 | 321072 | 10.277 | pbv # | | 91 |
| 20) Carbon Disulfide | 6.341 | 76 | 1113637 | 10.209 | ppbv | | 100 |
| 21) Trichlorotrifluoroethane | 6.275 | 101 | 927111 | 10.363 | ppbv | | 99 |
| 22) Trans-1,2-Dichloroethene | 6.740 | 61 | 527031 | 10.535 | ppbv | | 96 |
| 23) 1,1-Dichloroethane | 6.871 | 63 | 664247 | 10.392 | ppbv | | 99 |
| 24) Methyl tert-butyl ethe... | 6.912 | 73 | 993805 | 10.816 | ppbv | | 99 |
| 25) Vinyl Acetate | 6.942 | 43 | 1489318 | 12.394 | ppbv# | | 99 |
| 26) Methyl Ethyl Ketone | 7.125 | 43 | 572709 | 10.368 | ppbv | | 97 |
| 27) Cis-1,2-Dichloroethene | 7.410 | 61 | 494727 | 10.899 | ppbv | | 96 |
| 28) Hexane | 7.529 | 57 | 523731 | 10.910 | ppbv | | 97 |
| 29) Chloroform | 7.592 | 83 | 862932 | 10.566 | ppbv | | 100 |
| 30) Ethyl acetate | 7.521 | 61 | 125593 | 8.404 | ppbv | | 99 |
| 31) Tetrahydrofuran | 7.870 | 42 | 326848 | 11.908 | ppbv | | 94 |
| 32) 1,2-Dichloroethane | 8.068 | 62 | 471215 | 10.577 | ppbv | | 98 |
| 33) 1,1,1-Trichloroethane | 8.227 | 97 | 857606 | 10.413 | ppbv | | 99 |
| 34) Benzene | 8.505 | 78 | 1242586 | 10.605 | ppbv | | 99 |
| 35) Carbon Tetrachloride | 8.600 | 117 | 960160 | 10.501 | ppbv | | 99 |
| 36) Cyclohexane | 8.678 | 84 | 484270 | 9.886 | ppbv | | 96 |
| 38) 1,2-dichloropropane | 8.990 | 63 | 476960 | 10.613 | ppbv | | 97 |
| 39) Bromodichloromethane | 9.102 | 83 | 940064 | 10.388 | ppbv | | 100 |
| 40) Trichloroethene | 9.119 | 130 | 682250 | 10.478 | ppbv | | 100 |
| 41) 2,2,4-trimethylpentane | 9.119 | 57 | 1651707 | 10.715 | ppbv | | 99 |
| 42) 1,4-Dioxane | 9.119 | 88 | 276285 | 10.706 | ppbv | | 94 |
| 43) Methyl methacrylate | 9.189 | 69 | 458296 | 10.866 | ppbv | | 98 |
| 44) Heptane | 9.249 | 43 | 554702 | 10.792 | ppbv | | 93 |

Data Path : H:\AIR2023\CHEM39\05MAY\23\
 Data File : 0523_20.D
 Acq On : 24 May 2023 1:15 am
 Operator :
 Sample : 10ppb lcs ; to1503t
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: May 25 16:52:43 2023
 Quant Method : H:\AIR2023\CHEM39\Methods\39_AIR_0523.M
 Quant Title : VOA Standards for 5 point calibration
 QLast Update : Thu May 25 16:51:55 2023
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------------------|--------|------|----------|--------|-------|----------|
| 45) cis-1,3-Dichloropropene | 9.603 | 75 | 743800 | 10.916 | ppbv | 99 |
| 46) 4-Methyl-2-pentanone(M... | 9.611 | 43 | 743306 | 11.010 | ppbv | 98 |
| 47) trans-1,3-Dichloropropene | 9.892 | 75 | 673465 | 11.326 | ppbv | 100 |
| 48) 1,1,2-Trichloroethane | 10.005 | 97 | 576608 | 10.532 | ppbv | 99 |
| 49) Toluene | 10.167 | 91 | 1659853 | 10.905 | ppbv | 99 |
| 50) Dibromochloromethane | 10.420 | 129 | 1186168 | 10.637 | ppbv | 99 |
| 51) 2-Hexanone(MBK) | 10.286 | 43 | 719113 | 11.304 | ppbv | 98 |
| 52) 1,2-Dibromoethane(EDB) | 10.561 | 107 | 974923 | 10.714 | ppbv | 100 |
| 53) Tetrachloroethene | 10.800 | 166 | 958979 | 10.652 | ppbv | 99 |
| 55) 1,1,1,2-Tetrachloroethane | 11.170 | 131 | 728633 | 9.795 | ppbv | 100 |
| 56) Chlorobenzene | 11.186 | 112 | 1390186 | 10.583 | ppbv | 98 |
| 57) Ethylbenzene | 11.368 | 91 | 2108661 | 10.922 | ppbv | 100 |
| 58) m,p-Xylene | 11.466 | 91 | 3325176 | 23.263 | ppbv | 99 |
| 59) Bromoform | 11.550 | 173 | 1352467 | 11.081 | ppbv | 100 |
| 60) Styrene | 11.671 | 104 | 1326983 | 11.678 | ppbv | 98 |
| 61) 1,1,2,2-Tetrachloroethane | 11.732 | 83 | 1270065 | 11.037 | ppbv | 99 |
| 62) o-Xylene | 11.732 | 91 | 1698047 | 11.349 | ppbv | 100 |
| 64) 2-Chlorotoluene | 12.338 | 126 | 612723 | 9.941 | ppbv | 99 |
| 65) Isopropylbenzene | 12.050 | 105 | 2236666 | 9.837 | ppbv | 99 |
| 66) n-Propylbenzene | 12.338 | 120 | 656392 | 10.084 | ppbv | 99 |
| 67) 4-Ethyltoluene | 12.422 | 105 | 2635400 | 11.898 | ppbv | 97 |
| 68) 1,3,5-Trimethylbenzene | 12.460 | 105 | 1879491 | 10.488 | ppbv | 94 |
| 69) 1,2,4-Trimethylbenzene | 12.711 | 105 | 2047631 | 11.520 | ppbv | 100 |
| 70) tert-butylbenzene | 12.711 | 119 | 2088750 | 10.145 | ppbv | 100 |
| 71) Benzyl chloride | 12.817 | 91 | 1649611 | 11.443 | ppbv | 98 |
| 72) 1,3-Dichlorobenzene | 12.832 | 146 | 1633299 | 11.680 | ppbv | 99 |
| 73) 1,4-Dichlorobenzene | 12.870 | 146 | 1405532 | 10.425 | ppbv | 99 |
| 74) sec-Butylbenzene | 12.878 | 105 | 2763634 | 9.927 | ppbv | 100 |
| 75) 4-Isopropyltoluene | 12.961 | 119 | 2533768 | 9.877 | ppbv | 100 |
| 76) 1,2-Dichlorobenzene | 13.090 | 146 | 1396963 | 11.016 | ppbv | 99 |
| 77) n-Butylbenzene | 13.227 | 91 | 2173540 | 9.925 | ppbv | 98 |
| 78) 1,2,4-Trichlorobenzene | 14.283 | 180 | 966625 | 9.415 | ppbv | 99 |
| 79) Naphthalene | 14.389 | 128 | 1961258 | 9.903 | ppbv# | 100 |
| 80) Hexachlorobutadiene | 14.640 | 225 | 1014362 | 9.432 | ppbv | 100 |
| 82] 1,2-Dichlorotetrafluor... | 4.349 | 85 | 1109384 | 10.426 | ppbv | 98 |
| 83] Vinyl Chloride(sim) | 4.465 | 62 | 372248 | 10.761 | ppbv | 100 |
| 84] Bromomethane(sim) | 4.828 | 94 | 423424 | 10.238 | ppbv | 99 |
| 85] Trichlorofluoromethane... | 5.568 | 101 | 1136146 | 10.265 | ppbv# | 100 |
| 86] 1,2-Dichloroethane(sim) | 8.068 | 62 | 471215 | 10.407 | ppbv | 98 |
| 87] 1,1,1-Trichloroethane(... | 8.230 | 97 | 924744 | 10.560 | ppbv# | 99 |
| 88] Benzene(sim) | 8.505 | 78 | 1242893 | 9.840 | ppbv | 99 |
| 89] Carbon Tetrachloride(sim) | 8.603 | 117 | 958659 | 10.553 | ppbv | 100 |
| 90] 1,1-Dichloroethene(sim) | 6.025 | 61 | 564995 | 10.219 | ppbv | 94 |
| 91] Trichlorotrifluoroetha... | 6.278 | 101 | 918166 | 10.154 | ppbv# | 99 |
| 92] Trans-1,2-Dichloroethe... | 6.740 | 61 | 527031 | 10.459 | ppbv | 96 |
| 93] 1,1-Dichloroethane(sim) | 6.874 | 63 | 699804 | 10.302 | ppbv | 100 |
| 94] Cis-1,2-Dichloroethene... | 7.410 | 61 | 494727 | 10.689 | ppbv | 96 |
| 95] Chloroform(sim) | 7.595 | 83 | 887203 | 9.706 | ppbv | 99 |
| 97] 1,2-dichloropropane(sim) | 8.984 | 63 | 495072 | 10.091 | ppbv | 97 |
| 98] Bromodichloromethane(sim) | 9.102 | 83 | 940064 | 9.776 | ppbv | 99 |
| 99] Trichloroethene(sim) | 9.122 | 130 | 676247 | 10.075 | ppbv | 99 |
| 100] 1,4-Dioxane(sim) | 9.119 | 88 | 276285 | 11.295 | ppbv | 94 |
| 101] cis-1,3-Dichloropropen... | 9.599 | 75 | 778909 | 11.190 | ppbv | 100 |
| 102] 1,1,2-Trichloroethane(... | 10.005 | 97 | 576608 | 9.878 | ppbv | 99 |
| 103] Dibromochloromethane(sim) | 10.423 | 129 | 1170412 | 10.399 | ppbv | 100 |

Data Path : H:\AIR2023\CHEM39\05MAY\23\
 Data File : 0523_20.D
 Acq On : 24 May 2023 1:15 am
 Operator :
 Sample : 10ppb lcs ; to1503t
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

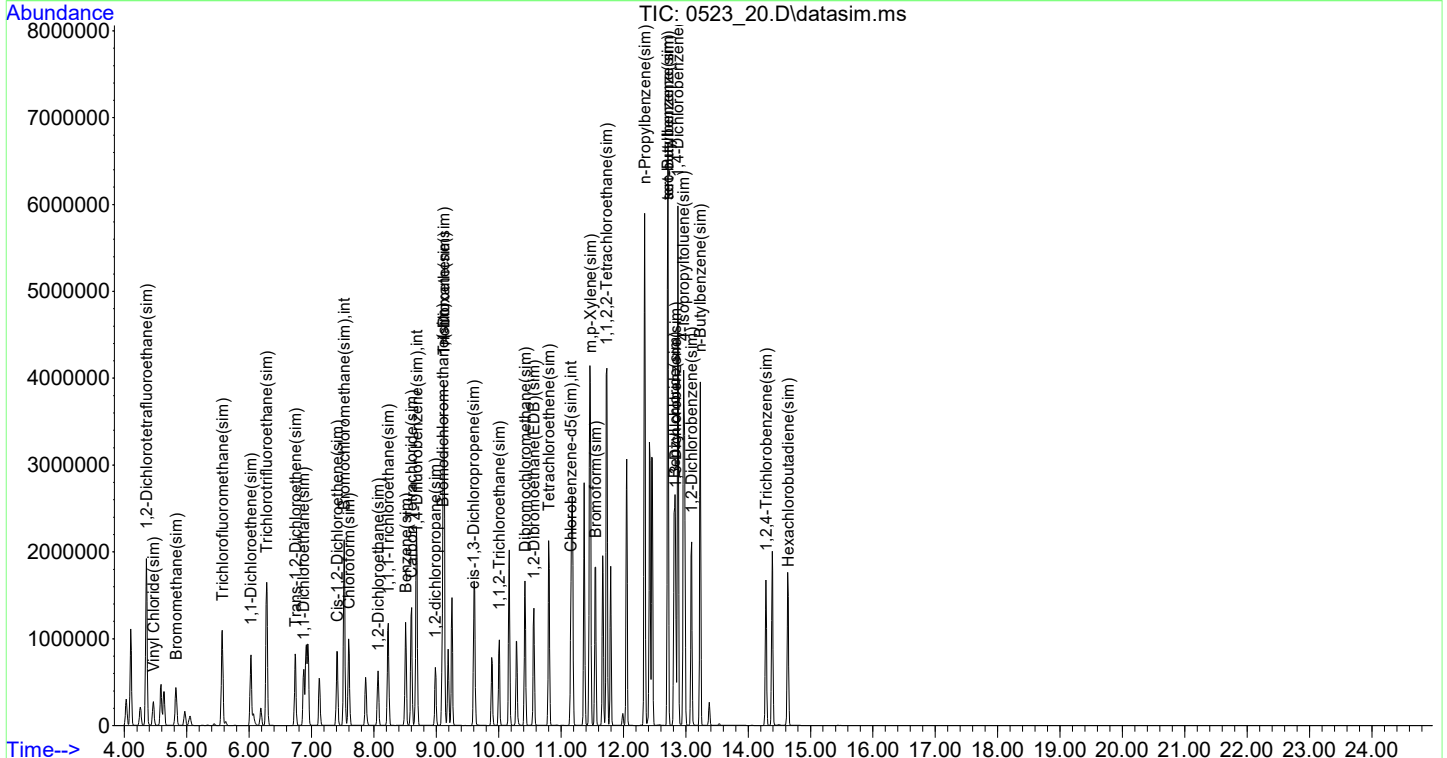
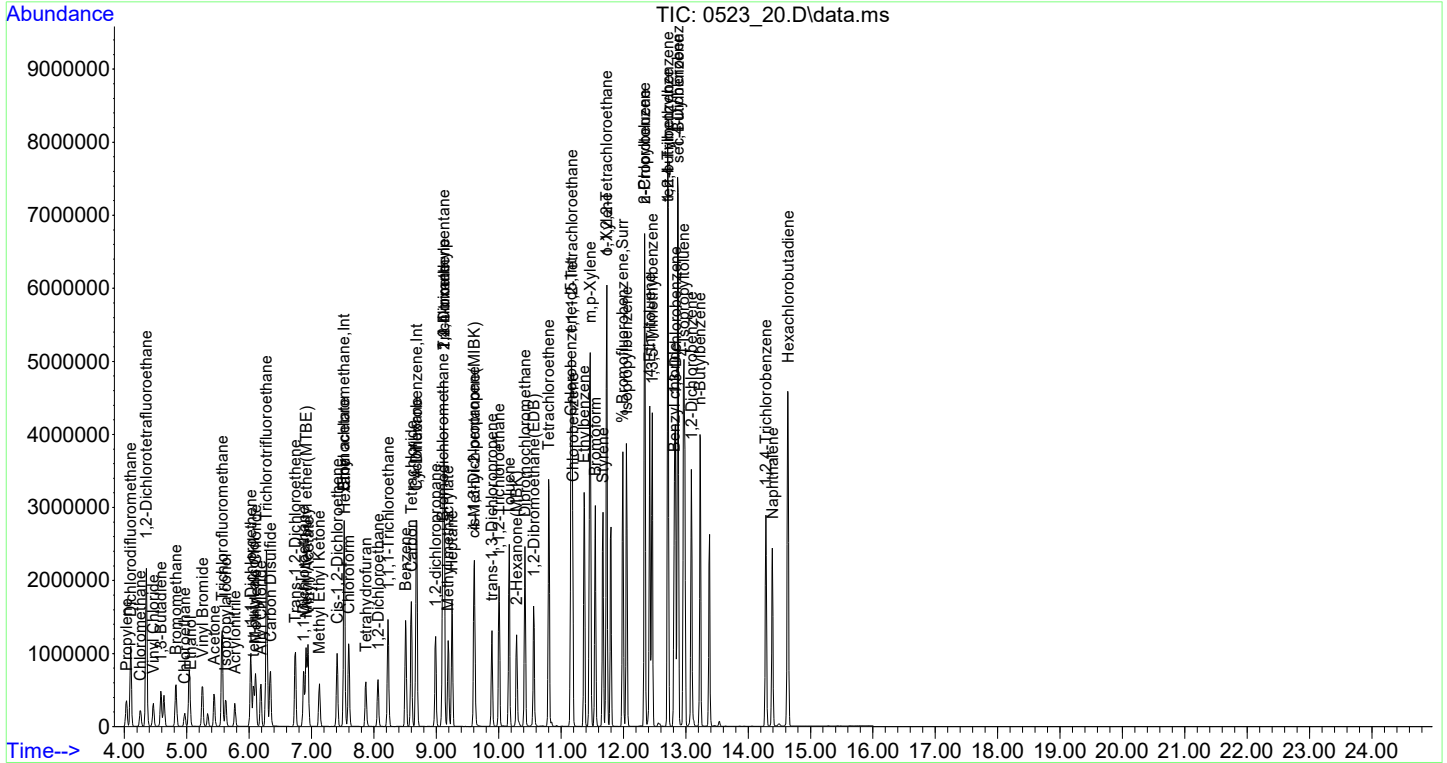
Quant Time: May 25 16:52:43 2023
 Quant Method : H:\AIR2023\CHEM39\Methods\39_AIR_0523.M
 Quant Title : VOA Standards for 5 point calibration
 QLast Update : Thu May 25 16:51:55 2023
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------------------|--------|------|----------|--------|-------|----------|
| 104] 1,2-Dibromoethane(EDB)... | 10.561 | 107 | 974923 | 10.529 | ppbv | 100 |
| 105] Tetrachloroethene(sim) | 10.803 | 166 | 940321 | 10.026 | ppbv | 99 |
| 107] Bromoform(sim) | 11.552 | 173 | 1286730 | 10.329 | ppbv | 100 |
| 108] m,p-Xylene(sim) | 11.466 | 91 | 3333180 | 22.946 | ppbv | 100 |
| 109] 1,1,2,2-Tetrachloroeth... | 11.727 | 83 | 1279548 | 10.105 | ppbv | 100 |
| 110] n-Propylbenzene(sim) | 12.338 | 120 | 656392 | 9.575 | ppbv | 99 |
| 111] tert-butylbenzene(sim) | 12.714 | 119 | 2276904 | 10.228 | ppbv# | 99 |
| 112] Benzyl chloride(sim) | 12.817 | 91 | 1648409 | 12.737 | ppbv | 98 |
| 113] 1,3-Dichlorobenzene(sim) | 12.835 | 146 | 1618113 | 10.700 | ppbv | 99 |
| 114] 1,4-Dichlorobenzene(sim) | 12.870 | 146 | 1405288 | 9.995 | ppbv | 99 |
| 115] sec-Butylbenzene(sim) | 12.714 | 105 | 2211211 | 12.194 | ppbv | 99 |
| 116] 4-Isopropyltoluene(sim) | 12.961 | 119 | 2533768 | 10.471 | ppbv | 100 |
| 117] 1,2-Dichlorobenzene(sim) | 13.093 | 146 | 1468145 | 11.058 | ppbv | 99 |
| 118] n-Butylbenzene(sim) | 13.227 | 91 | 2173540 | 10.898 | ppbv | 98 |
| 119] 1,2,4-Trichlorobenzene... | 14.286 | 180 | 1001451 | 8.838 | ppbv | 100 |
| 120] Hexachlorobutadiene(sim) | 14.640 | 225 | 1014287 | 8.146 | ppbv | 100 |

(#)out of range (m)manual integration reviewed by analyst (+)signals summed

Data Path : H:\AIR2023\CHEM39\05MAY\23\
 Data File : 0523_20.D
 Acq On : 24 May 2023 1:15 am
 Operator :
 Sample : 10ppb lcs ; to1503t
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: May 25 16:52:43 2023
 Quant Method : H:\AIR2023\CHEM39\Methods\39_AIR_0523.M
 Quant Title : VOA Standards for 5 point calibration
 QLast Update : Thu May 25 16:51:55 2023
 Response via : Initial Calibration



ATTACHMENT D

ANNOTATED SUMMARY FORMS

1
AIR ANALYSIS DATA SHEET

CLIENT ID

IA-4

| | | | |
|--------------|----------|------------------|-------------------|
| Client: | ACT | Lab: | Phoenix Env. Labs |
| SDG No.: | GCO73252 | Lab Sample ID: | CO73252 |
| Canister: | 28570 | Lab File ID: | 0811_13.D |
| Instrument: | CHEM39 | Column: | TX-1 ; #10157 |
| Purge Volume | 200 (cc) | Date Received: | 08/11/23 |
| Matrix: | AIR | Date Analyzed: | 08/11/23 |
| | | Dilution Factor: | 1 |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|-------------|-------------------------------|-------|----|-------|-------|---|
| 115-07-1 | Propylene | 0.581 | U | 0.581 | 0.581 | r |
| 75-71-8 | Dichlorodifluoromethane | 0.471 | | 0.202 | 0.202 | r |
| 74-87-3 | Chloromethane | 0.485 | U | 0.485 | 0.485 | r |
| 106-99-0 | 1,3-Butadiene | 0.452 | U | 0.452 | 0.452 | r |
| 75-00-3 | Chloroethane | 0.379 | U | 0.379 | 0.379 | r |
| 64-17-5 | Ethanol | 9.67 | J- | 0.531 | 0.531 | r |
| 67-64-1 | Acetone | 18.6 | | 0.421 | 0.421 | r |
| 75-69-4 | Trichlorofluoromethane | 0.293 | | 0.178 | 0.178 | r |
| 67-63-0 | Isopropylalcohol | 1.89 | S | 0.407 | 0.407 | r |
| 107-13-1 | Acrylonitrile | 0.461 | U | 0.461 | 0.461 | r |
| 75-09-2 | Methylene Chloride | 0.863 | U | 0.863 | 0.863 | r |
| 75-15-0 | Carbon Disulfide | 0.321 | U | 0.321 | 0.321 | r |
| 1634-04-4 | Methyl tert-butyl ether(MTBE) | 0.278 | U | 0.278 | 0.278 | r |
| 78-93-3 | Methyl Ethyl Ketone | 2.60 | | 0.339 | 0.339 | r |
| 156-59-2 | Cis-1,2-Dichloroethene | 7.78 | | 0.051 | 0.051 | r |
| 110-54-3 | Hexane | 1.74 | S | 0.284 | 0.284 | r |
| 67-66-3 | Chloroform | 0.605 | | 0.205 | 0.205 | r |
| 141-78-6 | Ethyl acetate | 0.278 | U | 0.278 | 0.278 | r |
| 109-99-9 | Tetrahydrofuran | 27.1 | | 0.339 | 0.339 | r |
| 71-43-2 | Benzene | 0.784 | | 0.313 | 0.313 | r |
| 110-82-7 | Cyclohexane | 0.291 | U | 0.291 | 0.291 | r |
| 79-01-6 | Trichloroethene | 1.33 | | 0.037 | 0.037 | r |
| 142-82-5 | Heptane | 0.397 | | 0.244 | 0.244 | r |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | 0.244 | U | 0.244 | 0.244 | r |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.221 | U | 0.221 | 0.221 | r |
| 108-88-3 | Toluene | 3.66 | | 0.266 | 0.266 | r |
| 591-78-6 | 2-Hexanone(MBK) | 0.244 | U | 0.244 | 0.244 | r |
| 127-18-4 | Tetrachloroethene | 2.61 | | 0.037 | 0.037 | r |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.146 | U | 0.146 | 0.146 | r |
| 108-90-7 | Chlorobenzene | 0.217 | U | 0.217 | 0.217 | r |
| 100-41-4 | Ethylbenzene | 0.531 | | 0.230 | 0.230 | r |
| 179601-23-1 | m,p-Xylene | 2.18 | | 0.230 | 0.230 | r |
| 100-42-5 | Styrene | 0.235 | U | 0.235 | 0.235 | r |
| 95-47-6 | o-Xylene | 0.862 | | 0.230 | 0.230 | r |
| 98-82-8 | Isopropylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 622-96-8 | 4-Ethyltoluene | 0.204 | U | 0.204 | 0.204 | r |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

1
AIR ANALYSIS DATA SHEET

CLIENT ID

IA-3

Client: ACT Lab: Phoenix Env. Labs

SDG No.: GCO73252 Lab Sample ID: CO73253

Canister: 23348 Lab File ID: 0816_08.D

Instrument: CHEM39 Column: TX-1 ; #10157 Date Received: 08/11/23

Purge Volume 200 (cc) Date Analyzed: 08/16/23

Matrix: AIR Dilution Factor: 1

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|-------------|-------------------------------|-------|---|-------|-------|---|
| 115-07-1 | Propylene | 0.581 | U | 0.581 | 0.581 | r |
| 75-71-8 | Dichlorodifluoromethane | 0.454 | | 0.202 | 0.202 | r |
| 74-87-3 | Chloromethane | 0.485 | U | 0.485 | 0.485 | r |
| 106-99-0 | 1,3-Butadiene | 0.452 | U | 0.452 | 0.452 | r |
| 75-00-3 | Chloroethane | 0.379 | U | 0.379 | 0.379 | r |
| 64-17-5 | Ethanol | 18.6 | | 0.531 | 0.531 | r |
| 67-64-1 | Acetone | 33.5 | | 0.421 | 0.421 | r |
| 75-69-4 | Trichlorofluoromethane | 0.315 | | 0.178 | 0.178 | r |
| 67-63-0 | Isopropylalcohol | 2.94 | | 0.407 | 0.407 | r |
| 107-13-1 | Acrylonitrile | 0.461 | U | 0.461 | 0.461 | r |
| 75-09-2 | Methylene Chloride | 0.863 | U | 0.863 | 0.863 | r |
| 75-15-0 | Carbon Disulfide | 0.321 | U | 0.321 | 0.321 | r |
| 1634-04-4 | Methyl tert-butyl ether(MTBE) | 0.278 | U | 0.278 | 0.278 | r |
| 78-93-3 | Methyl Ethyl Ketone | 6.58 | | 0.339 | 0.339 | r |
| 156-59-2 | Cis-1,2-Dichloroethene | 7.92 | | 0.051 | 0.051 | r |
| 110-54-3 | Hexane | 2.84 | | 0.284 | 0.284 | r |
| 67-66-3 | Chloroform | 0.515 | | 0.205 | 0.205 | r |
| 141-78-6 | Ethyl acetate | 0.278 | U | 0.278 | 0.278 | r |
| 109-99-9 | Tetrahydrofuran | 45.0 | J | 0.339 | 0.339 | |
| 71-43-2 | Benzene | 1.05 | | 0.313 | 0.313 | r |
| 110-82-7 | Cyclohexane | 0.291 | U | 0.291 | 0.291 | r |
| 79-01-6 | Trichloroethene | 1.42 | | 0.037 | 0.037 | r |
| 142-82-5 | Heptane | 0.693 | | 0.244 | 0.244 | r |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | 0.244 | U | 0.244 | 0.244 | r |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.221 | U | 0.221 | 0.221 | r |
| 108-88-3 | Toluene | 5.61 | | 0.266 | 0.266 | r |
| 591-78-6 | 2-Hexanone(MBK) | 0.244 | U | 0.244 | 0.244 | r |
| 127-18-4 | Tetrachloroethene | 2.61 | | 0.037 | 0.037 | r |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.146 | U | 0.146 | 0.146 | r |
| 108-90-7 | Chlorobenzene | 0.217 | U | 0.217 | 0.217 | r |
| 100-41-4 | Ethylbenzene | 0.833 | | 0.230 | 0.230 | r |
| 179601-23-1 | m,p-Xylene | 3.43 | | 0.230 | 0.230 | r |
| 100-42-5 | Styrene | 0.235 | U | 0.235 | 0.235 | r |
| 95-47-6 | o-Xylene | 1.32 | | 0.230 | 0.230 | r |
| 98-82-8 | Isopropylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 622-96-8 | 4-Ethyltoluene | 1.60 | | 0.204 | 0.204 | r |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

1
AIR ANALYSIS DATA SHEET

CLIENT ID

IA-3

| | | | |
|--------------|----------|------------------|-------------------|
| Client: | ACT | Lab: | Phoenix Env. Labs |
| SDG No.: | GCO73252 | Lab Sample ID: | CO73253 |
| Canister: | 23348 | Lab File ID: | 0816_08.D |
| Instrument: | CHEM39 | Column: | FX-1 ; #10157 |
| Purge Volume | 200 (cc) | Date Received: | 08/11/23 |
| Matrix: | AIR | Date Analyzed: | 08/16/23 |
| | | Dilution Factor: | 1 |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|------------|------------------------------------|-------|---|-------|-------|---|
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.619 | | 0.204 | 0.204 | r |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane(sim) | 0.143 | U | 0.143 | 0.143 | r |
| 75-01-4 | Vinyl Chloride(sim) | 0.089 | | 0.078 | 0.078 | r |
| 74-83-9 | Bromomethane(sim) | 0.258 | U | 0.258 | 0.258 | r |
| 107-06-2 | 1,2-Dichloroethane(sim) | 0.247 | U | 0.247 | 0.247 | r |
| 71-55-6 | 1,1,1-Trichloroethane(sim) | 0.183 | U | 0.183 | 0.183 | r |
| 56-23-5 | Carbon Tetrachloride(sim) | 0.080 | | 0.032 | 0.032 | r |
| 75-35-4 | 1,1-Dichloroethene(sim) | 0.051 | U | 0.051 | 0.051 | r |
| 76-13-1 | Trichlorotrifluoroethane(sim) | 0.131 | U | 0.131 | 0.131 | r |
| 156-60-5 | Trans-1,2-Dichloroethene(sim) | 0.252 | U | 0.252 | 0.252 | r |
| 75-34-3 | 1,1-Dichloroethane(sim) | 0.247 | U | 0.247 | 0.247 | r |
| 78-87-5 | 1,2-dichloropropane(sim) | 0.217 | U | 0.217 | 0.217 | r |
| 75-27-4 | Bromodichloromethane(sim) | 0.149 | U | 0.149 | 0.149 | r |
| 123-91-1 | 1,4-Dioxane(sim) | 0.278 | U | 0.278 | 0.278 | r |
| 10061-01-5 | cis-1,3-Dichloropropene(sim) | 0.221 | U | 0.221 | 0.221 | r |
| 79-00-5 | 1,1,2-Trichloroethane(sim) | 0.183 | U | 0.183 | 0.183 | r |
| 124-48-1 | Dibromochloromethane(sim) | 0.118 | U | 0.118 | 0.118 | r |
| 106-93-4 | 1,2-Dibromoethane(EDB)(sim) | 0.130 | U | 0.130 | 0.130 | r |
| 75-25-2 | Bromoform(sim) | 0.097 | U | 0.097 | 0.097 | r |
| 79-34-5 | 1,1,2,2-Tetrachloroethane(sim) | 0.146 | U | 0.146 | 0.146 | r |
| 100-44-7 | Benzyl chloride(sim) | 0.193 | U | 0.193 | 0.193 | r |
| 541-73-1 | 1,3-Dichlorobenzene(sim) | 0.166 | U | 0.166 | 0.166 | r |
| 106-46-7 | 1,4-Dichlorobenzene(sim) | 0.166 | U | 0.166 | 0.166 | r |
| 135-98-8 | sec-Butylbenzene(sim) | 0.182 | U | 0.182 | 0.182 | r |
| 99-87-6 | 4-Isopropyltoluene(sim) | 0.182 | U | 0.182 | 0.182 | r |
| 95-50-1 | 1,2-Dichlorobenzene(sim) | 0.166 | U | 0.166 | 0.166 | r |
| 104-51-8 | n-Butylbenzene(sim) | 0.182 | U | 0.182 | 0.182 | r |
| 120-82-1 | 1,2,4-Trichlorobenzene(sim) | 0.135 | U | 0.135 | 0.135 | r |
| 87-68-3 | Hexachlorobutadiene(sim) | 0.094 | U | 0.094 | 0.094 | r |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

JR 12/8/2025

1
AIR ANALYSIS DATA SHEET

CLIENT ID

IA-7

Client: ACT Lab: Phoenix Env. Labs

SDG No.: GCO73252 Lab Sample ID: CO73254

Canister: 489 Lab File ID: 0816_10.D

Instrument: CHEM39 Column: FX-1 ; #10157 Date Received: 08/11/23

Purge Volume 200 (cc) Date Analyzed: 08/16/23

Matrix: AIR Dilution Factor: 1

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|-------------|-------------------------------|-------|---|-------|-------|---|
| 115-07-1 | Propylene | 0.581 | U | 0.581 | 0.581 | r |
| 75-71-8 | Dichlorodifluoromethane | 0.449 | | 0.202 | 0.202 | r |
| 74-87-3 | Chloromethane | 0.485 | U | 0.485 | 0.485 | r |
| 106-99-0 | 1,3-Butadiene | 0.452 | U | 0.452 | 0.452 | r |
| 75-00-3 | Chloroethane | 0.379 | U | 0.379 | 0.379 | r |
| 64-17-5 | Ethanol | 36.6 | S | 0.531 | 0.531 | r |
| 67-64-1 | Acetone | 57.6 | J | 0.421 | 0.421 | |
| 75-69-4 | Trichlorofluoromethane | 0.413 | | 0.178 | 0.178 | r |
| 67-63-0 | Isopropylalcohol | 4.01 | | 0.407 | 0.407 | r |
| 107-13-1 | Acrylonitrile | 0.461 | U | 0.461 | 0.461 | r |
| 75-09-2 | Methylene Chloride | 0.863 | U | 0.863 | 0.863 | r |
| 75-15-0 | Carbon Disulfide | 0.321 | U | 0.321 | 0.321 | r |
| 1634-04-4 | Methyl tert-butyl ether(MTBE) | 0.278 | U | 0.278 | 0.278 | r |
| 78-93-3 | Methyl Ethyl Ketone | 9.58 | | 0.339 | 0.339 | r |
| 156-59-2 | Cis-1,2-Dichloroethene | 5.38 | | 0.051 | 0.051 | r |
| 110-54-3 | Hexane | 5.36 | | 0.284 | 0.284 | r |
| 67-66-3 | Chloroform | 0.457 | | 0.205 | 0.205 | r |
| 141-78-6 | Ethyl acetate | 0.278 | U | 0.278 | 0.278 | r |
| 109-99-9 | Tetrahydrofuran | 68.1 | J | 0.339 | 0.339 | |
| 71-43-2 | Benzene | 1.74 | | 0.313 | 0.313 | r |
| 110-82-7 | Cyclohexane | 0.291 | U | 0.291 | 0.291 | r |
| 79-01-6 | Trichloroethene | 1.02 | | 0.037 | 0.037 | r |
| 142-82-5 | Heptane | 1.63 | | 0.244 | 0.244 | r |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | 0.244 | U | 0.244 | 0.244 | r |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.221 | U | 0.221 | 0.221 | r |
| 108-88-3 | Toluene | 8.97 | | 0.266 | 0.266 | r |
| 591-78-6 | 2-Hexanone(MBK) | 0.244 | U | 0.244 | 0.244 | r |
| 127-18-4 | Tetrachloroethene | 2.03 | | 0.037 | 0.037 | r |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.146 | U | 0.146 | 0.146 | r |
| 108-90-7 | Chlorobenzene | 0.217 | U | 0.217 | 0.217 | r |
| 100-41-4 | Ethylbenzene | 1.39 | | 0.230 | 0.230 | r |
| 179601-23-1 | m,p-Xylene | 5.77 | | 0.230 | 0.230 | r |
| 100-42-5 | Styrene | 0.235 | U | 0.235 | 0.235 | r |
| 95-47-6 | o-Xylene | 2.19 | | 0.230 | 0.230 | r |
| 98-82-8 | Isopropylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 622-96-8 | 4-Ethyltoluene | 0.204 | U | 0.204 | 0.204 | r |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

December 8, 2025

Mr. Jason Stewart
Advanced Cleanup Technologies
228 Park Ave S PMB 34864
New York, New York 10003

Re: Data Usability Summary Report – Phoenix Environmental Laboratories, Inc. – GCP07983 V1

Dear Mr. Stewart:

The evaluation of analytical data by Phoenix Environmental Laboratories for project 9268-BXNY, which were reported in a single data package under Job No. GCP07983 V1 has been completed. The following samples were reported.

IA-17 OA-4 SS-1B

Analysis was performed in accordance with EPA Method TO-15 (volatile organics). The review was performed to the extent possible, in accordance with the analytical method, and “DER-10/ Technical Guidance for Site Investigation and Remediation”. Professional judgment is applied as necessary and appropriate. National Functional Guidelines for Organic Data Review was consulted as needed. Qualifiers consistent with those defined by EPA Region 2 are applied as necessary and appropriate.

Below is the Data Usability Summary Report (DUSR) associated with these samples.

| Data Usability Summary Report | |
|---|--------------------------------|
| 1. Is the data package complete as defined under the requirements for the most current DEC ASP Category B or USEPA CLP data deliverables? | No – see Documentation section |
| 2. Have all holding times been met? | Yes |
| 3. Do all the QC data; blanks, instrument tunings, calibration standards, calibration verifications, surrogate recoveries, spike recoveries, replicate analyses, laboratory controls and sample data fall within the protocol required limits and specifications? | No -see following sections |
| 4. Have all of the data been generated using established and agreed upon analytical protocols? | Yes |
| 5. Does an evaluation of the raw data confirm the results provided in the data summary sheet and the quality control verification forms? | Yes |
| 6. Have the correct data qualifiers been used and are they consistent with the most current DEC ASP? | Yes |

| | |
|---|-----|
| 7. Have any quality control (QC) exceedances been specifically noted in the DUSR and have the corresponding QC summary sheet from the data package been attached to the DUSR? | Yes |
|---|-----|

Overall Evaluation

Based on the data review effort, results are usable, with the following qualifications. For samples that are qualified as estimated (J-, UJ), detected results may be biased low. False negatives may exist in non-detect results. Sample results that are qualified as estimated (J+) may be biased high. For samples that are qualified as estimated with any combination of (J), (J-) and/or (J+), the (J) qualifier takes precedence and is applied to the sample result. It is not possible to determine the direction of the bias and the overall effect on the result.

- The results for carbon tetrachloride, trans-1,3-dichloropropene, dibromochloromethane, 1,1,1,2-tetrachloroethane and bromoform in all samples are qualified as estimated (J-, UJ) due to low response in the continuing calibration verification (CCV) standard.
- The result for ethanol in all samples is qualified as estimated (J-, UJ) due to low recovery in the laboratory control sample (LCS).
- The result for ethanol in IA-17 is qualified as estimated (J) because the reported concentration is outside the calibration range.
- The results for acetone in all samples are qualified as estimated (J+) because the sample concentration is less than ten times the acetone concentration in the clean canister concentration.

Qualifier definitions are provided in Attachment A. A copy of the chain of custody record is provided in Attachment B. Pages from the data package illustrating the exceedances and issues described in this validation report are included in Attachment C. Annotated Summary Forms are included in Attachment D detailing qualifications resulting from the data review effort.

The following components were reviewed, where applicable:

- Chain of Custody
- Receiving conditions
- Holding times
- Preservation
- Analyte lists
- Reporting limits
- Requested methods
- Units, and
- Sample related quality control data:
 - Method, instrument blanks
 - Clean canister certification
 - Field blanks

- Surrogate recoveries
- LCS recoveries
- Internal standard area response
- Duplicates
- Analyte Identification
- Instrument related quality control data:
 - Instrument tunes
 - Calibration summaries

The following sections of the report detail only quality control exceedances that impacted results. Where a quality control item exceeded control limits but there is no impact to the samples results, these are not detailed in the report.

Documentation: A completeness review of the data package was performed, and the data package was determined to be a complete Category B data package, with the following exception:

- The data package does not include summary forms and raw data for the second source initial calibration verification (ICV) standard. The laboratory was contacted and provided the ICV data and is included in attachment C.

The following documentation issues were observed during the review:

- Results for ethanol, acetone, isopropyl alcohol, and hexane in all samples are flagged 'S' which the data deliverable identifies this flag as 'This compound is a solvent that is used in the laboratory. Laboratory contamination is suspected if concentration is less than five times the reporting level.' Where the sample detections are greater than five times the reporting limit, the validator removed the 'S' qualifier.

Holding Times, Preservation, Sample Integrity:

A copy of the applicable chain of custody (COC) record was included in the data package, documenting sample collection date of September 21, 2023. The samples were received on September 25, 2023. Samples were received intact and analyzed within the method specified hold time.

A. Volatile Organics

1. Calibration

One initial calibration (IC) was performed in support of the sample analysis. All relative response factors (RRFs) and relative standard deviations (RSDs) or correlation coefficients are acceptable. A second source ICV standard was analyzed after the IC, and all percent differences are acceptable (70-

130%R). One continuing calibration was analyzed, and percent differences (%D) are acceptable, with exceptions noted below.

| Analyte | %R | Associated Sample | Qualifier Applied |
|-----------------------------------|------|-------------------|-------------------|
| <i>CCV CHEM20 9/25/2025 11:10</i> | | | |
| Carbon Tetrachloride | 35.0 | IA-17 | J-, UJ |
| trans-1,3-Dichloropropene | 31.9 | OA-4 | |
| Dibromochloromethane | 33.3 | SS-1B | |
| 1,1,1,2-Tetrachloroethane | 33.5 | | |
| Bromoform | 35.3 | | |

In all instances, the percent difference represents a decrease in instrument sensitivity. The results for carbon tetrachloride, trans-1,3-dichloropropene, dibromochloromethane, 1,1,1,2-tetrachloroethane and bromoform in all samples are qualified as estimated (J-, UJ) due to low response in the CCV.

2. Laboratory Control Sample (LCS)

One LCS was prepared and analyzed with the sample. The LCS is evaluated using control limit of 70-130%R. All recoveries are acceptable, with the following exception:

| Analyte | LCS %R | Affected Sample | Qualifier Applied |
|--------------------|--------|------------------------|-------------------|
| <i>CO71323 LCS</i> | | | |
| Ethanol | 66 | IA-17 OA-4 SS-1B | J-, UJ |

*The summary form provided by the laboratory reports LCS recoveries based on the full scan analysis. The validator reviewed the raw data to determine the recovery in the SIM analysis and recoveries are acceptable.

The result for ethanol in all samples is qualified as estimated (J-, UJ) due to low recovery in the LCS.

3. Compound Quantitation

The result for ethanol in IA-17 is qualified as estimated (J) because the reported concentration is outside the calibration range. Results are quantitatively uncertain.

4. Clean Canister Certification

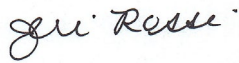
Acetone is detected in two of the clean canister certification batches as presented below.

| Batch Canister | Analyte | Conc. (ppbv) |
|-----------------------|----------------|---------------------|
| Batch #3402 | Acetone | 0.613 |
| Batch #3043 | Acetone | 0.464 |

The data package does not indicate which canisters are associated with which batch canister certification data. Based on professional judgement, the highest acetone value (0.613 ppbv) is used to evaluate sample results. On this basis, the results for acetone in all samples are qualified as estimated (J+) because the sample concentration is less than ten times the clean canister concentration.

No other sample results are qualified. Please feel free to contact me at (908) 370-3431 or richjერიrossi513@gmail.com if you have any questions regarding this data package review report or need further information.

Sincerely,



Jeri L Rossi, CEAC

Environmental Consulting Chemist

ATTACHMENT A

Qualifier Definitions

EPA Qualifier Definitions

- U The analyte was analyzed for but was not detected above the level of the reported sample quantitation limit.
- J The result is an estimated quantity. The associated numerical value is the approximate concentration of the analyte in the sample.
- J+ The result is an estimated quantity, but the result may be biased high.
- J- The result is an estimated quantity, but the result may be biased low.
- NJ The analyte has been “tentatively identified” or “presumptively” as present and the associated numerical value is the estimated concentration in the sample.
- UJ The analyte was analyzed for but was not detected. The reported quantitation limit is approximate and may be inaccurate or imprecise.
- R The data are unusable. The sample results are rejected due to serious deficiencies in meeting QC criteria. The analyte may or may not be present in the sample.

ATTACHMENT B

CHAIN OF CUSTODY (COC)



597 5th Middle Turnpike, P.O. Box 370, Manhattan, CT 06340
Telephone: 860-645-1102 • Fax: 860-645-0521

ACT

CHAIN OF CUSTODY RECORD

AIR ANALYSES

860-645-1102

email: greg@phoenixlabs.com

P.O. #

Page 1 of 1

Data Delivery:

Fax #:

Email:

rayb@act-earth or Jason@act-earth

Phone #:

| | | |
|---|----------------------------|---|
| Report to: Joseph Squeglia | Project Name: 9628 - BXNY | Data Format: (Circle) Equis Excel Other: |
| Customer: Advanced Cleanup Technologies, Inc. | Invoice to: Karen Friedman | Requested Deliverable: RCP ASP CAT B |
| Address: 228 Park Ave S PMB 34864 | | MCP NJ Deliverables |
| 15001 New York, New York 10003 | Sampled by: Yunmee Han | Quote Number: |

| Phoenix ID # | Client Sample ID | Canister ID # | Canister Size (L) | Outgoing Canister Pressure ("Hg) | Incoming Canister Pressure ("Hg) | Flow Regulator ID # | Flow Controller Setting (mL/min) | Sampling Start Time | Sampling End Time | Sample Start Date | Canister Pressure at Start ("Hg) | Canister Pressure at End ("Hg) | Ambient/Indoor Air | Soil Gas | Grab (G) Composite (C) | TO-15 | APH |
|-------------------------------|--------------------------------|------------------|-------------------|----------------------------------|----------------------------------|---------------------|----------------------------------|---------------------|-------------------|-------------------|----------------------------------|--------------------------------|--------------------|----------|------------------------|-------|-----|
| THIS SECTION FOR LAB USE ONLY | | | | | | | | | | | | | | | | | |
| | | 23348 | 6.0L | -30 | | | 7013 | 3.13 | | | | | | | | | |
| | | 489 | 6.0L | -30 | | | 10559 | 3.93 | | | | | | | | | |
| 07983 | 1A-17 | 226 | 6.0L | -30 | -8 | 10651 | 3.09 | 1402 | | 9/21/23 | -30 | -9 | ✓ | | | | ✓ |
| 07984 | 0A-4 | 28588 | 6.0L | -30 | -8 | 6994 | 3.08 | 1410 | | 9/21/23 | -30 | -9 | ✓ | | | | ✓ |
| | Extra canister, do not analyze | 21357 | 6.0L | -30 | | 3252 | 3.3 | | | | | | | | | | |
| 07985 | SS-18 | 28568 | 6.0L | -30 | -8 | 3221 | 3.17 | 1358 | | 9/21/23 | -30 | -10 | ✓ | | | | ✓ |
| | | 227 | 6.0L | -30 | | | 10605 | 3.3 | | | | | | | | | |

| | | | | |
|-----------------------------------|-----------------|---------------|------------|---|
| Relinquished by: Yunmee Han Zm | Accepted by: Zm | Date: 9-25-23 | Time: 1033 | I attest that all media released by Phoenix Environmental Laboratories, Inc. have been received in good working condition and agree to the terms and conditions as listed on the back of this document. |
| | | Date: 9-25-23 | Time: 1719 | |

| | | | | | | | |
|--------------------------------|--|--|--|--|---------------------------------------|---|---|
| State Where Samples Collected: | Turnaround Time: 1 Day <input type="checkbox"/> 2 Day* <input type="checkbox"/> 3 Day* <input type="checkbox"/> 4 Day* <input type="checkbox"/> 5 Day* <input type="checkbox"/> Standard <input checked="" type="checkbox"/> | Requested Criteria: (Please Circle) CT: TAC I/C TAC RES SVVC I/C SVVC RES GWV I/C GWV CES | MA: Indoor Air Residential Ind/Commercial Soil Gas: Residential Ind/Commercial | NI: Indoor Air Residential Ind/Commercial | NY: Vapor Intrusion NYS DEC Eds | PA: Indoor Air Residential Non-residential | VT: Indoor Air Residential Industrial Sub-slab Residential Industrial |
|--------------------------------|--|--|--|--|---------------------------------------|---|---|

SPECIAL INSTRUCTIONS, OC REQUIREMENTS, REGULATORY INFORMATION:
(7) - 6.0L 24 hr
*NYS DEC Eds
* Received & Analyzed (EP)

*SURCHARGES MAY APPLY

01/14/2024

Phoenix Environmental Laboratories, Inc.

Page 22 of 211

ATTACHMENT C

**SELECTED PAGES FROM DATA PACKAGE –
QC EXCEEDANCES AND VALIDATION ISSUES**

3
AIR LCS RECOVERY

Lab Name: Phoenix Environmental Labs Client: ACT

Lab Code: Phoenix Case No: _____ SAS No: _____ SDG No GCP07983

LCS - Client Id: CP07053 LCS

| COMPOUND | SPIKE ADDED (ppbv) | | LCS CONCENTRATION (ppbv) | LCS % REC # | QC. LIMITS REC. | |
|-------------------------------|--------------------|--|--------------------------|-------------|-----------------|-----|
| Propylene | 10 | | 8.667 | 87 | 70 | 130 |
| Dichlorodifluoromethane | 10 | | 9.126 | 91 | 70 | 130 |
| Chloromethane | 10 | | 8.373 | 84 | 70 | 130 |
| 1,2-Dichlorotetrafluoroethane | 10 | | 9.270 | 93 | 70 | 130 |
| Vinyl Chloride | 10 | | 8.611 | 86 | 70 | 130 |
| 1,3-Butadiene | 10 | | 8.354 | 84 | 70 | 130 |
| Bromomethane | 10 | | 9.214 | 92 | 70 | 130 |
| Chloroethane | 10 | | 8.429 | 84 | 70 | 130 |
| Ethanol | 10 | | 13.60 | 136 * | 70 | 130 |
| Acetone | 10 | | 8.286 | 83 | 70 | 130 |
| Trichlorofluoromethane | 10 | | 9.792 | 98 | 70 | 130 |
| Isopropylalcohol | 10 | | 8.467 | 85 | 70 | 130 |
| Acrylonitrile | 10 | | 9.065 | 91 | 70 | 130 |
| 1,1-Dichloroethene | 10 | | 8.773 | 88 | 70 | 130 |
| Methylene Chloride | 10 | | 8.519 | 85 | 70 | 130 |
| Carbon Disulfide | 10 | | 8.858 | 89 | 70 | 130 |
| Trichlorotrifluoroethane | 10 | | 9.297 | 93 | 70 | 130 |
| Trans-1,2-Dichloroethene | 10 | | 8.783 | 88 | 70 | 130 |
| 1,1-Dichloroethane | 10 | | 8.769 | 88 | 70 | 130 |
| Methyl tert-butyl ether(MTBE) | 10 | | 9.246 | 92 | 70 | 130 |
| Methyl Ethyl Ketone | 10 | | 9.598 | 96 | 70 | 130 |
| Cis-1,2-Dichloroethene | 10 | | 8.624 | 86 | 70 | 130 |
| Hexane | 10 | | 8.907 | 89 | 70 | 130 |
| Chloroform | 10 | | 9.819 | 98 | 70 | 130 |
| Ethyl acetate | 10 | | 6.063 | 61 * | 70 | 130 |
| Tetrahydrofuran | 10 | | 9.652 | 97 | 70 | 130 |
| 1,2-Dichloroethane | 10 | | 8.975 | 90 | 70 | 130 |
| 1,1,1-Trichloroethane | 10 | | 8.311 | 83 | 70 | 130 |
| Benzene | 10 | | 9.279 | 93 | 70 | 130 |
| Carbon Tetrachloride | 10 | | 7.015 | 70 | 70 | 130 |
| Cyclohexane | 10 | | 9.309 | 93 | 70 | 130 |
| 1,2-dichloropropane | 10 | | 9.508 | 95 | 70 | 130 |
| Bromodichloromethane | 10 | | 9.272 | 93 | 70 | 130 |
| Trichloroethene | 10 | | 10.28 | 103 | 70 | 130 |
| 1,4-Dioxane | 10 | | 9.830 | 98 | 70 | 130 |
| Heptane | 10 | | 9.000 | 90 | 70 | 130 |
| cis-1,3-Dichloropropene | 10 | | 9.138 | 91 | 70 | 130 |
| 4-Methyl-2-pentanone(MIBK) | 10 | | 9.658 | 97 | 70 | 130 |
| trans-1,3-Dichloropropene | 10 | | 9.191 | 92 | 70 | 130 |
| 1,1,2-Trichloroethane | 10 | | 10.21 | 102 | 70 | 130 |
| Toluene | 10 | | 9.828 | 98 | 70 | 130 |
| Dibromochloromethane | 10 | | 8.856 | 89 | 70 | 130 |
| 2-Hexanone(MBK) | 10 | | 9.962 | 100 | 70 | 130 |
| 1,2-Dibromoethane(EDB) | 10 | | 10.03 | 100 | 70 | 130 |

FORM III AIR

Response Factor Report CHEM20

Method Path : H:\AIR2023\CHEM20\Methods\
 Method File : 20_AIR_0921.M
 Title : VOA Standards for 5 point calibration

| Method | Path | File | Title | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 | |
|--------|------|------|-----------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|----|-------|---------|-------|
| 97) | | | 1,2-dichloropr... | 0.582 | 0.546 | 0.567 | 0.528 | 0.495 | 0.466 | 0.477 | 0.583 | | | 0.505 | 0.528 | 8.45 |
| 98) | | | Bromdichlorom... | 0.706 | 0.790 | 0.750 | 0.642 | 0.607 | 0.557 | 0.576 | 0.693 | | | 0.758 | 0.676 | 12.45 |
| 99) | | | Trichloroethen... | 0.465 | 0.473 | 0.476 | 0.436 | 0.418 | 0.397 | 0.399 | 0.483 | | | 0.509 | 0.451 | 8.84 |
| 100) | | | 1,4-Dioxane(sim) | | | 0.249 | 0.186 | 0.192 | 0.177 | 0.165 | 0.217 | | | | 0.198 | 15.37 |
| 101) | | | cis-1,3-Dichlo... | 0.567 | 0.548 | 0.586 | 0.550 | 0.531 | 0.501 | 0.517 | | | | 0.565 | 0.546 | 5.14 |
| 102) | | | 1,1,2-Trichlor... | 0.475 | 0.407 | 0.433 | 0.350 | 0.328 | 0.311 | 0.331 | 0.412 | | | 0.443 | 0.388 | 15.18 |
| 103) | | | Dibromchlorom... | 0.640 | 0.610 | 0.676 | 0.611 | 0.607 | 0.585 | 0.610 | | | | 0.614 | 0.619 | 4.44 |
| 104) | | | 1,2-Dibrommeth... | 0.668 | 0.684 | 0.683 | 0.594 | 0.577 | 0.557 | 0.549 | 0.692 | | | 0.618 | 0.625 | 9.28 |
| 105) | | | Tetrachloroeth... | 0.689 | 0.651 | 0.664 | 0.613 | 0.579 | 0.538 | 0.545 | 0.682 | 0.619 | | 0.762 | 0.634 | 10.98 |
| 106) | int | | Chlorobenzene-d5(sim) | -----ISTD----- | | | | | | | | | | | | |
| 107) | | | Bromform(sim) | 1.122 | 1.046 | 1.158 | 1.044 | 1.014 | 0.927 | 0.983 | | | | 1.251 | 1.068 | 9.71 |
| 108) | | | m p-Xylene(sim) | 2.334 | 2.175 | 2.395 | 2.057 | 2.016 | 1.787 | 1.835 | 2.336 | | | 2.451 | 2.154 | 11.35 |
| 109) | | | 1,1,2,2-Tetrac... | 1.617 | 1.609 | 1.681 | 1.560 | 1.477 | 1.167 | 1.159 | 1.804 | | | 1.838 | 1.546 | 15.80 |
| 112) | | | Benzyl chlorid... | 1.754 | 1.790 | 1.910 | 1.664 | 1.659 | 1.181 | 1.101 | | | | | 1.580 | 19.78 |
| 113) | | | 1,3-Dichlorobe... | 1.958 | 1.805 | 1.930 | 1.728 | 1.660 | 1.178 | 1.101 | | | | | 1.623 | 21.38 |
| 114) | qfi | | 1,4-Dichlorobe... | 0.049 | 0.075 | 0.118 | 0.202 | 0.523 | 0.830 | 2.894 | | | | | Coef R2 | 0.99 |
| 115) | | | sec-Butylbenze... | 4.392 | 4.184 | 4.509 | 4.146 | 3.945 | 3.384 | 3.404 | 4.331 | | | 5.122 | 4.158 | 13.04 |
| 116) | | | 4-Isopropyltol... | 4.097 | 4.156 | 4.261 | 3.828 | 3.527 | 2.955 | 2.994 | | | | 4.974 | 3.849 | 17.64 |
| 117) | | | 1,2-Dichlorobe... | 1.963 | 1.894 | 1.870 | 1.730 | 1.638 | 1.028 | 0.949 | 1.910 | | | 2.544 | 1.725 | 28.31 |
| 118) | | | n-Butylbenzene... | 3.890 | 3.961 | 3.659 | 3.274 | 3.125 | 2.416 | 2.453 | | | | 4.623 | 3.425 | 22.26 |
| 119) | qfi | | 1,2,4-Trichlor... | 0.053 | 0.062 | 0.110 | 0.205 | 0.516 | 0.878 | 2.616 | | | | | Coef R2 | 1.00 |
| 120) | qfi | | Hexachlorobuta... | 0.054 | 0.066 | 0.126 | 0.211 | 0.537 | 0.749 | 2.878 | | | | 0.038 | Coef R2 | 0.98@ |

(#, \$, @)=Out of Range l=linear lf=linear(0,0) q=Quadratic qf=Quadratic(0,0)

7A
AIR CONTINUING CALIBRATION CHECK

Lab Name: Phoenix Environmental Labs Client: _____
 Lab Code: Phoenix Case No.: _____ SAS No.: _____ SDG No.: GCP07983
 Instrument: CHEM20 Calibration Date: 09/25/23 Time: 11:10
 Lab File Id: 0925_02.D Init. Calib. Date(s): 09/20/23 09/20/23
 Heated Purge (Y/N): Y Init. Calib. Times: 14:15 20:09
 GC Column: RTX-1 60M Method File: 20_AIR_0921.M

| COMPOUND | RRF | RRF1 | RRF MIN | %D | % D LIMITS |
|-------------------------------|-------|-------|---------|--------|------------|
| Propylene | 1.302 | 1.263 | | 3.0 | 30 |
| Dichlorodifluoromethane | 2.693 | 2.676 | | 0.6 | 30 |
| Chloromethane | 1.566 | 1.497 | | 4.4 | 30 |
| 1,2-Dichlorotetrafluoroethane | 2.636 | 2.708 | | -2.7 | 30 |
| Vinyl Chloride | 1.432 | 1.389 | | 3.0 | 30 |
| 1,3-Butadiene | 1.374 | 1.306 | | 4.9 | 30 |
| Bromomethane | 0.971 | 1.008 | | -3.8 | 30 |
| Chloroethane | 0.628 | 0.588 | | 6.4 | 30 |
| Ethanol | 0.424 | 0.417 | | 1.7 | 30 |
| Acetone | 2.689 | 2.671 | | 0.7 | 30 |
| Trichlorofluoromethane | 2.875 | 2.897 | | -0.8 | 30 |
| Isopropylalcohol | 3.310 | 3.235 | | 2.3 | 30 |
| Acrylonitrile | 1.557 | 1.329 | | 14.6 | 30 |
| 1,1-Dichloroethene | 2.527 | 2.471 | | 2.2 | 30 |
| Methylene Chloride | 2.105 | 2.083 | | 1.0 | 30 |
| Carbon Disulfide | 2.862 | 2.920 | | -2.0 | 30 |
| Trichlorotrifluoroethane | 2.182 | 2.270 | | -4.0 | 30 |
| Trans-1,2-Dichloroethene | 2.346 | 2.288 | | 2.5 | 30 |
| 1,1-Dichloroethane | 2.631 | 2.594 | | 1.4 | 30 |
| Methyl tert-butyl ether(MTBE) | 2.862 | 2.891 | | -1.0 | 30 |
| Methyl Ethyl Ketone | 3.598 | 3.394 | | 5.7 | 30 |
| Cis-1,2-Dichloroethene | 2.382 | 2.191 | | 8.0 | 30 |
| Hexane | 2.781 | 2.595 | | 6.7 | 30 |
| Chloroform | 2.296 | 2.483 | | -8.1 | 30 |
| Ethyl acetate | 0.768 | 0.635 | | 17.3 | 30 |
| Tetrahydrofuran | 1.958 | 2.015 | | -2.9 | 30 |
| 1,2-Dichloroethane | 2.168 | 2.153 | | 0.7 | 30 |
| 1,1,1-Trichloroethane | 2.410 | 2.026 | | 15.9 | 30 |
| Benzene | 3.362 | 3.413 | | -1.5 | 30 |
| Carbon Tetrachloride | 2.552 | 1.660 | | 35.0 # | 30 |
| Cyclohexane | 1.324 | 1.439 | | -8.7 | 30 |
| 1,2-dichloropropane | 0.473 | 0.438 | | 7.4 | 30 |
| Bromodichloromethane | 0.627 | 0.501 | | 20.1 | 30 |
| Trichloroethene | 0.394 | 0.406 | | -3.0 | 30 |
| 1,4-Dioxane | 0.188 | 0.191 | | -1.6 | 30 |

(*) Recommended RRF not met (+) %D exceeds criteria % (#) %D exceeds (maximum) criteria
 %D: 20% of target compounds are allowed to be above criteria %, but must be less than the (maximum) %D
 (#) Maximum %D not met.

7B
AIR CONTINUING CALIBRATION CHECK

Lab Name: Phoenix Environmental Labs Client: _____
 Lab Code: Phoenix Case No.: _____ SAS No.: _____ SDG No.: GCP07983
 Instrument: CHEM20 Calibration Date: 09/25/23 Time: 11:10
 Lab File Id: 0925_02.D Init. Calib. Date(s): 09/20/23 09/20/23
 Heated Purge (Y/N): Y Init. Calib. Times: 14:15 20:09
 GC Column: RTX-1 60M Method File: 20_AIR_0921.M

| COMPOUND | RRF | RRF1 | RRF MIN | %D | % D LIMITS |
|------------------------------------|-------|-------|---------|--------|------------|
| Heptane | 0.926 | 0.896 | | 3.2 | 30 |
| cis-1,3-Dichloropropene | 0.515 | 0.399 | | 22.5 | 30 |
| 4-Methyl-2-pentanone(MIBK) | 1.143 | 1.121 | | 1.9 | 30 |
| trans-1,3-Dichloropropene | 0.474 | 0.323 | | 31.9 # | 30 |
| 1,1,2-Trichloroethane | 0.354 | 0.355 | | -0.3 | 30 |
| Toluene | 1.092 | 1.072 | | 1.8 | 30 |
| Dibromochloromethane | 0.634 | 0.423 | | 33.3 # | 30 |
| 2-Hexanone(MBK) | 1.067 | 0.980 | | 8.2 | 30 |
| 1,2-Dibromoethane(EDB) | 0.604 | 0.508 | | 15.9 | 30 |
| Tetrachloroethene | 0.523 | 0.529 | | -1.1 | 30 |
| 1,1,1,2-Tetrachloroethane | 0.832 | 0.553 | | 33.5 # | 30 |
| Chlorobenzene | 1.499 | 1.615 | | -7.7 | 30 |
| Ethylbenzene | 2.578 | 2.598 | | -0.8 | 30 |
| m,p-Xylene | 1.916 | 1.643 | | 14.2 | 30 |
| Bromoform | 0.998 | 0.646 | | 35.3 # | 30 |
| Styrene | 1.478 | 1.527 | | -3.3 | 30 |
| 1,1,2,2-Tetrachloroethane | 1.337 | 1.386 | | -3.7 | 30 |
| o-Xylene | 2.047 | 2.100 | | -2.6 | 30 |
| Isopropylbenzene | 2.837 | 2.997 | | -5.6 | 30 |
| 4-Ethyltoluene | 2.749 | 2.845 | | -3.5 | 30 |
| 1,3,5-Trimethylbenzene | 2.241 | 2.450 | | -9.3 | 30 |
| 1,2,4-Trimethylbenzene | 2.275 | 2.342 | | -2.9 | 30 |
| Benzyl chloride | 1.714 | 1.396 | | 18.6 | 30 |
| 1,3-Dichlorobenzene | 1.410 | 1.566 | | -11.1 | 30 |
| 1,4-Dichlorobenzene | 1.373 | 1.480 | | -7.8 | 30 |
| sec-Butylbenzene | 3.547 | 3.660 | | -3.2 | 30 |
| 4-Isopropyltoluene | 3.336 | 3.581 | | -7.3 | 30 |
| 1,2-Dichlorobenzene | 1.336 | 1.424 | | -6.6 | 30 |
| n-Butylbenzene | 2.860 | 2.975 | | -4.0 | 30 |
| 1,2,4-Trichlorobenzene | 1.133 | 1.109 | | 2.1 | 30 |
| Hexachlorobutadiene | 1.235 | 1.180 | | 4.5 | 30 |
| 1,2-Dichlorotetrafluoroethane(sim) | 2.468 | 2.569 | | -4.1 | 30 |
| Vinyl Chloride(sim) | 1.425 | 1.433 | | -0.6 | 30 |
| Bromomethane(sim) | 0.919 | 0.957 | | -4.1 | 30 |
| Trichlorofluoromethane(sim) | 2.749 | 2.848 | | -3.6 | 30 |

(*) Recommended RRF not met (+) %D exceeds criteria % (#) %D exceeds (maximum) criteria
 %D: 20% of target compounds are allowed to be above criteria %, but must be less than the (maximum) %D
 (#) Maximum %D not met.

7B
AIR CONTINUING CALIBRATION CHECK

Lab Name: Phoenix Environmental Labs Client: _____
 Lab Code: Phoenix Case No.: _____ SAS No.: _____ SDG No.: GCP07983
 Instrument: CHEM20 Calibration Date: 09/25/23 Time: 11:10
 Lab File Id: 0925_02.D Init. Calib. Date(s): 09/20/23 09/20/23
 Heated Purge (Y/N): Y Init. Calib. Times: 14:15 20:09
 GC Column: RTX-1 60M Method File: 20_AIR_0921.M

| COMPOUND | RRF | RRF1 | RRF MIN | %D | % D LIMITS |
|---------------------------------|-------|-------|---------|--------|------------|
| 1,2-Dichloroethane(sim) | 1.962 | 2.043 | | -4.1 | 30 |
| 1,1,1-Trichloroethane(sim) | 2.420 | 2.042 | | 15.6 | 30 |
| Benzene(sim) | 3.378 | 3.239 | | 4.1 | 30 |
| Carbon Tetrachloride(sim) | 2.350 | 1.649 | | 29.8 | 30 |
| 1,1-Dichloroethene(sim) | 2.401 | 2.344 | | 2.4 | 30 |
| Trichlorotrifluoroethane(sim) | 2.190 | 2.287 | | -4.4 | 30 |
| Trans-1,2-Dichloroethene(sim) | 2.215 | 2.170 | | 2.0 | 30 |
| 1,1-Dichloroethane(sim) | 2.691 | 2.696 | | -0.2 | 30 |
| Cis-1,2-Dichloroethene(sim) | 2.217 | 2.079 | | 6.2 | 30 |
| Chloroform(sim) | 2.288 | 2.476 | | -8.2 | 30 |
| 1,2-dichloropropane(sim) | 0.528 | 0.473 | | 10.4 | 30 |
| Bromodichloromethane(sim) | 0.676 | 0.501 | | 25.9 | 30 |
| Trichloroethene(sim) | 0.451 | 0.439 | | 2.7 | 30 |
| 1,4-Dioxane(sim) | 0.198 | 0.191 | | 3.5 | 30 |
| cis-1,3-Dichloropropene(sim) | 0.546 | 0.442 | | 19.0 | 30 |
| 1,1,2-Trichloroethane(sim) | 0.388 | 0.355 | | 8.5 | 30 |
| Dibromochloromethane(sim) | 0.619 | 0.452 | | 27.0 | 30 |
| 1,2-Dibromoethane(EDB)(sim) | 0.625 | 0.508 | | 18.7 | 30 |
| Tetrachloroethene(sim) | 0.634 | 0.609 | | 3.9 | 30 |
| Bromoform(sim) | 1.068 | 0.755 | | 29.3 | 30 |
| m,p-Xylene(sim) | 2.154 | 2.088 | | 3.1 | 30 |
| 1,1,2,2-Tetrachloroethane(sim) | 1.546 | 1.465 | | 5.2 | 30 |
| Benzyl chloride(sim) | 1.580 | 1.500 | | 5.1 | 30 |
| 1,3-Dichlorobenzene(sim) | 1.623 | 1.771 | | -9.1 | 30 |
| 1,4-Dichlorobenzene(sim) qfi | 1.000 | 1.18 | | -18.0 | 20 |
| sec-Butylbenzene(sim) | 4.158 | 3.914 | | 5.9 | 30 |
| 4-Isopropyltoluene(sim) | 3.849 | 3.586 | | 6.8 | 30 |
| 1,2-Dichlorobenzene(sim) | 1.725 | 1.536 | | 11.0 | 30 |
| n-Butylbenzene(sim) | 3.425 | 2.988 | | 12.8 | 30 |
| 1,2,4-Trichlorobenzene(sim) qfi | 1.000 | 0.82 | | 18.0 | 20 |
| Hexachlorobutadiene(sim) qfi | 1.000 | 0.75 | | 25.0 # | 20 |
| % Bromofluorobenzene | 1.327 | 1.383 | | -4.2 | 30 |
| | | | | | |
| | | | | | |
| | | | | | |

(*) Recommended RRF not met (+) %D exceeds criteria % (#) %D exceeds (maximum) criteria
 %D: 20% of target compounds are allowed to be above criteria %, but must be less than the (maximum) %D
 (#) Maximum %D not met.

7A
AIR CONTINUING CALIBRATION CHECK

Lab Name: Phoenix Environmental Labs Client: _____
 Lab Code: Phoenix Case No.: _____ SAS No.: _____ SDG No.: GCP07983
 Instrument: CHEM20 Calibration Date: 09/25/23 Time: 11:10
 Lab File Id: 0925_02.D Init. Calib. Date(s): 09/20/23 09/20/23
 Heated Purge (Y/N): Y Init. Calib. Times: 14:15 20:09
 GC Column: RTX-1 60M Method File: 20_AIR_0921.M

| COMPOUND | RRF | RRF1 | RRF MIN | %D | % D LIMITS |
|-------------------------------|-------|-------|---------|--------|------------|
| Propylene | 1.302 | 1.263 | | 3.0 | 30 |
| Dichlorodifluoromethane | 2.693 | 2.676 | | 0.6 | 30 |
| Chloromethane | 1.566 | 1.497 | | 4.4 | 30 |
| 1,2-Dichlorotetrafluoroethane | 2.636 | 2.708 | | -2.7 | 30 |
| Vinyl Chloride | 1.432 | 1.389 | | 3.0 | 30 |
| 1,3-Butadiene | 1.374 | 1.306 | | 4.9 | 30 |
| Bromomethane | 0.971 | 1.008 | | -3.8 | 30 |
| Chloroethane | 0.628 | 0.588 | | 6.4 | 30 |
| Ethanol | 0.424 | 0.417 | | 1.7 | 30 |
| Acetone | 2.689 | 2.671 | | 0.7 | 30 |
| Trichlorofluoromethane | 2.875 | 2.897 | | -0.8 | 30 |
| Isopropylalcohol | 3.310 | 3.235 | | 2.3 | 30 |
| Acrylonitrile | 1.557 | 1.329 | | 14.6 | 30 |
| 1,1-Dichloroethene | 2.527 | 2.471 | | 2.2 | 30 |
| Methylene Chloride | 2.105 | 2.083 | | 1.0 | 30 |
| Carbon Disulfide | 2.862 | 2.920 | | -2.0 | 30 |
| Trichlorotrifluoroethane | 2.182 | 2.270 | | -4.0 | 30 |
| Trans-1,2-Dichloroethene | 2.346 | 2.288 | | 2.5 | 30 |
| 1,1-Dichloroethane | 2.631 | 2.594 | | 1.4 | 30 |
| Methyl tert-butyl ether(MTBE) | 2.862 | 2.891 | | -1.0 | 30 |
| Methyl Ethyl Ketone | 3.598 | 3.394 | | 5.7 | 30 |
| Cis-1,2-Dichloroethene | 2.382 | 2.191 | | 8.0 | 30 |
| Hexane | 2.781 | 2.595 | | 6.7 | 30 |
| Chloroform | 2.296 | 2.483 | | -8.1 | 30 |
| Ethyl acetate | 0.768 | 0.635 | | 17.3 | 30 |
| Tetrahydrofuran | 1.958 | 2.015 | | -2.9 | 30 |
| 1,2-Dichloroethane | 2.168 | 2.153 | | 0.7 | 30 |
| 1,1,1-Trichloroethane | 2.410 | 2.026 | | 15.9 | 30 |
| Benzene | 3.362 | 3.413 | | -1.5 | 30 |
| Carbon Tetrachloride | 2.552 | 1.660 | | 35.0 # | 30 |
| Cyclohexane | 1.324 | 1.439 | | -8.7 | 30 |
| 1,2-dichloropropane | 0.473 | 0.438 | | 7.4 | 30 |
| Bromodichloromethane | 0.627 | 0.501 | | 20.1 | 30 |
| Trichloroethene | 0.394 | 0.406 | | -3.0 | 30 |
| 1,4-Dioxane | 0.188 | 0.191 | | -1.6 | 30 |

(*) Recommended RRF not met (+) %D exceeds criteria % (#) %D exceeds (maximum) criteria
 %D: 20% of target compounds are allowed to be above criteria %, but must be less than the (maximum) %D
 (#) Maximum %D not met.

7B
AIR CONTINUING CALIBRATION CHECK

Lab Name: Phoenix Environmental Labs Client: _____

Lab Code: Phoenix Case No.: _____ SAS No.: _____ SDG No.: GCP07983

Instrument: CHEM20 Calibration Date: 09/25/23 Time: 11:10

Lab File Id: 0925_02.D Init. Calib. Date(s): 09/20/23 09/20/23

Heated Purge (Y/N): Y Init. Calib. Times: 14:15 20:09

GC Column: RTX-1 60M Method File: 20_AIR_0921.M

| COMPOUND | RRF | RRF1 | RRF MIN | %D | % D LIMITS |
|------------------------------------|-------|-------|---------|--------|------------|
| Heptane | 0.926 | 0.896 | | 3.2 | 30 |
| cis-1,3-Dichloropropene | 0.515 | 0.399 | | 22.5 | 30 |
| 4-Methyl-2-pentanone(MIBK) | 1.143 | 1.121 | | 1.9 | 30 |
| trans-1,3-Dichloropropene | 0.474 | 0.323 | | 31.9 # | 30 |
| 1,1,2-Trichloroethane | 0.354 | 0.355 | | -0.3 | 30 |
| Toluene | 1.092 | 1.072 | | 1.8 | 30 |
| Dibromochloromethane | 0.634 | 0.423 | | 33.3 # | 30 |
| 2-Hexanone(MBK) | 1.067 | 0.980 | | 8.2 | 30 |
| 1,2-Dibromoethane(EDB) | 0.604 | 0.508 | | 15.9 | 30 |
| Tetrachloroethene | 0.523 | 0.529 | | -1.1 | 30 |
| 1,1,1,2-Tetrachloroethane | 0.832 | 0.553 | | 33.5 # | 30 |
| Chlorobenzene | 1.499 | 1.615 | | -7.7 | 30 |
| Ethylbenzene | 2.578 | 2.598 | | -0.8 | 30 |
| m,p-Xylene | 1.916 | 1.643 | | 14.2 | 30 |
| Bromoform | 0.998 | 0.646 | | 35.3 # | 30 |
| Styrene | 1.478 | 1.527 | | -3.3 | 30 |
| 1,1,2,2-Tetrachloroethane | 1.337 | 1.386 | | -3.7 | 30 |
| o-Xylene | 2.047 | 2.100 | | -2.6 | 30 |
| Isopropylbenzene | 2.837 | 2.997 | | -5.6 | 30 |
| 4-Ethyltoluene | 2.749 | 2.845 | | -3.5 | 30 |
| 1,3,5-Trimethylbenzene | 2.241 | 2.450 | | -9.3 | 30 |
| 1,2,4-Trimethylbenzene | 2.275 | 2.342 | | -2.9 | 30 |
| Benzyl chloride | 1.714 | 1.396 | | 18.6 | 30 |
| 1,3-Dichlorobenzene | 1.410 | 1.566 | | -11.1 | 30 |
| 1,4-Dichlorobenzene | 1.373 | 1.480 | | -7.8 | 30 |
| sec-Butylbenzene | 3.547 | 3.660 | | -3.2 | 30 |
| 4-Isopropyltoluene | 3.336 | 3.581 | | -7.3 | 30 |
| 1,2-Dichlorobenzene | 1.336 | 1.424 | | -6.6 | 30 |
| n-Butylbenzene | 2.860 | 2.975 | | -4.0 | 30 |
| 1,2,4-Trichlorobenzene | 1.133 | 1.109 | | 2.1 | 30 |
| Hexachlorobutadiene | 1.235 | 1.180 | | 4.5 | 30 |
| 1,2-Dichlorotetrafluoroethane(sim) | 2.468 | 2.569 | | -4.1 | 30 |
| Vinyl Chloride(sim) | 1.425 | 1.433 | | -0.6 | 30 |
| Bromomethane(sim) | 0.919 | 0.957 | | -4.1 | 30 |
| Trichlorofluoromethane(sim) | 2.749 | 2.848 | | -3.6 | 30 |

(*) Recommended RRF not met (+) %D exceeds criteria % (#) %D exceeds (maximum) criteria
%D: 20% of target compounds are allowed to be above criteria %, but must be less than the (maximum) %D
(#) Maximum %D not met.

7B
AIR CONTINUING CALIBRATION CHECK

Lab Name: Phoenix Environmental Labs Client: _____
 Lab Code: Phoenix Case No.: _____ SAS No.: _____ SDG No.: GCP07983
 Instrument: CHEM20 Calibration Date: 09/25/23 Time: 11:10
 Lab File Id: 0925_02.D Init. Calib. Date(s): 09/20/23 09/20/23
 Heated Purge (Y/N): Y Init. Calib. Times: 14:15 20:09
 GC Column: RTX-1 60M Method File: 20_AIR_0921.M

| COMPOUND | RRF | RRF1 | RRF MIN | %D | % D LIMITS |
|---------------------------------|-------|-------|---------|--------|------------|
| 1,2-Dichloroethane(sim) | 1.962 | 2.043 | | -4.1 | 30 |
| 1,1,1-Trichloroethane(sim) | 2.420 | 2.042 | | 15.6 | 30 |
| Benzene(sim) | 3.378 | 3.239 | | 4.1 | 30 |
| Carbon Tetrachloride(sim) | 2.350 | 1.649 | | 29.8 | 30 |
| 1,1-Dichloroethene(sim) | 2.401 | 2.344 | | 2.4 | 30 |
| Trichlorotrifluoroethane(sim) | 2.190 | 2.287 | | -4.4 | 30 |
| Trans-1,2-Dichloroethene(sim) | 2.215 | 2.170 | | 2.0 | 30 |
| 1,1-Dichloroethane(sim) | 2.691 | 2.696 | | -0.2 | 30 |
| Cis-1,2-Dichloroethene(sim) | 2.217 | 2.079 | | 6.2 | 30 |
| Chloroform(sim) | 2.288 | 2.476 | | -8.2 | 30 |
| 1,2-dichloropropane(sim) | 0.528 | 0.473 | | 10.4 | 30 |
| Bromodichloromethane(sim) | 0.676 | 0.501 | | 25.9 | 30 |
| Trichloroethene(sim) | 0.451 | 0.439 | | 2.7 | 30 |
| 1,4-Dioxane(sim) | 0.198 | 0.191 | | 3.5 | 30 |
| cis-1,3-Dichloropropene(sim) | 0.546 | 0.442 | | 19.0 | 30 |
| 1,1,2-Trichloroethane(sim) | 0.388 | 0.355 | | 8.5 | 30 |
| Dibromochloromethane(sim) | 0.619 | 0.452 | | 27.0 | 30 |
| 1,2-Dibromoethane(EDB)(sim) | 0.625 | 0.508 | | 18.7 | 30 |
| Tetrachloroethene(sim) | 0.634 | 0.609 | | 3.9 | 30 |
| Bromoform(sim) | 1.068 | 0.755 | | 29.3 | 30 |
| m,p-Xylene(sim) | 2.154 | 2.088 | | 3.1 | 30 |
| 1,1,2,2-Tetrachloroethane(sim) | 1.546 | 1.465 | | 5.2 | 30 |
| Benzyl chloride(sim) | 1.580 | 1.500 | | 5.1 | 30 |
| 1,3-Dichlorobenzene(sim) | 1.623 | 1.771 | | -9.1 | 30 |
| 1,4-Dichlorobenzene(sim) qfi | 1.000 | 1.18 | | -18.0 | 20 |
| sec-Butylbenzene(sim) | 4.158 | 3.914 | | 5.9 | 30 |
| 4-Isopropyltoluene(sim) | 3.849 | 3.586 | | 6.8 | 30 |
| 1,2-Dichlorobenzene(sim) | 1.725 | 1.536 | | 11.0 | 30 |
| n-Butylbenzene(sim) | 3.425 | 2.988 | | 12.8 | 30 |
| 1,2,4-Trichlorobenzene(sim) qfi | 1.000 | 0.82 | | 18.0 | 20 |
| Hexachlorobutadiene(sim) qfi | 1.000 | 0.75 | | 25.0 # | 20 |
| % Bromofluorobenzene | 1.327 | 1.383 | | -4.2 | 30 |
| | | | | | |
| | | | | | |
| | | | | | |

(*) Recommended RRF not met (+) %D exceeds criteria % (#) %D exceeds (maximum) criteria
 %D: 20% of target compounds are allowed to be above criteria %, but must be less than the (maximum) %D
 (#) Maximum %D not met.

ATTACHMENT D

ANNOTATED SUMMARY FORMS

1
AIR ANALYSIS DATA SHEET

CLIENT ID

IA-17

| | | | |
|--------------|----------|------------------|-------------------|
| Client: | ACT | Lab: | Phoenix Env. Labs |
| SDG No.: | GCP07983 | Lab Sample ID: | CP07983 |
| Canister: | 226 | Lab File ID: | 0925_24.D |
| Instrument: | CHEM20 | Column: | RTX-1 60M |
| | | Date Received: | 09/25/23 |
| Purge Volume | 200 (cc) | Date Analyzed: | 09/25/23 |
| Matrix: | AIR | Dilution Factor: | 1 |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|-------------|------------------------------------|-------|----|-------|-------|---|
| 115-07-1 | Propylene | 0.581 | U | 0.581 | 0.581 | r |
| 75-71-8 | Dichlorodifluoromethane | 0.524 | | 0.202 | 0.202 | r |
| 74-87-3 | Chloromethane | 0.526 | | 0.485 | 0.485 | r |
| 106-99-0 | 1,3-Butadiene | 0.452 | U | 0.452 | 0.452 | r |
| 75-00-3 | Chloroethane | 0.379 | U | 0.379 | 0.379 | r |
| 64-17-5 | Ethanol | 43.1 | J | 0.531 | 0.531 | r |
| 67-64-1 | Acetone | 5.65 | J+ | 0.421 | 0.421 | r |
| 75-69-4 | Trichlorofluoromethane | 0.239 | | 0.178 | 0.178 | r |
| 67-63-0 | Isopropylalcohol | 1.25 | S | 0.407 | 0.407 | r |
| 107-13-1 | Acrylonitrile | 0.461 | U | 0.461 | 0.461 | r |
| 75-09-2 | Methylene Chloride | 0.863 | U | 0.863 | 0.863 | r |
| 75-15-0 | Carbon Disulfide | 0.321 | U | 0.321 | 0.321 | r |
| 1634-04-4 | Methyl tert-butyl ether(MTBE) | 0.278 | U | 0.278 | 0.278 | r |
| 78-93-3 | Methyl Ethyl Ketone | 0.339 | U | 0.339 | 0.339 | r |
| 110-54-3 | Hexane | 0.284 | U | 0.284 | 0.284 | r |
| 141-78-6 | Ethyl acetate | 0.278 | U | 0.278 | 0.278 | r |
| 109-99-9 | Tetrahydrofuran | 0.339 | U | 0.339 | 0.339 | r |
| 71-43-2 | Benzene | 0.313 | U | 0.313 | 0.313 | r |
| 110-82-7 | Cyclohexane | 0.291 | U | 0.291 | 0.291 | r |
| 142-82-5 | Heptane | 0.244 | U | 0.244 | 0.244 | r |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | 0.244 | U | 0.244 | 0.244 | r |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.221 | U | 0.221 | 0.221 | r |
| 108-88-3 | Toluene | 0.623 | | 0.266 | 0.266 | r |
| 591-78-6 | 2-Hexanone(MBK) | 0.244 | U | 0.244 | 0.244 | r |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.146 | UJ | 0.146 | 0.146 | r |
| 108-90-7 | Chlorobenzene | 0.217 | U | 0.217 | 0.217 | r |
| 100-41-4 | Ethylbenzene | 0.230 | U | 0.230 | 0.230 | r |
| 179601-23-1 | m,p-Xylene | 0.443 | | 0.230 | 0.230 | r |
| 100-42-5 | Styrene | 0.235 | U | 0.235 | 0.235 | r |
| 95-47-6 | o-Xylene | 0.230 | U | 0.230 | 0.230 | r |
| 98-82-8 | Isopropylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 622-96-8 | 4-Ethyltoluene | 0.204 | U | 0.204 | 0.204 | r |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane(sim) | 0.143 | U | 0.143 | 0.143 | r |
| 75-01-4 | Vinyl Chloride(sim) | 0.078 | U | 0.078 | 0.078 | r |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

1
AIR ANALYSIS DATA SHEET

CLIENT ID

IA-17

| | | | |
|--------------|----------|------------------|-------------------|
| Client: | ACT | Lab: | Phoenix Env. Labs |
| SDG No.: | GCP07983 | Lab Sample ID: | CP07983 |
| Canister: | 226 | Lab File ID: | 0925_24.D |
| Instrument: | CHEM20 | Column: | RTX-1 60M |
| Purge Volume | 200 (cc) | Date Received: | 09/25/23 |
| Matrix: | AIR | Date Analyzed: | 09/25/23 |
| | | Dilution Factor: | 1 |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|------------|--------------------------------|-------|----|-------|-------|---|
| 74-83-9 | Bromomethane(sim) | 0.258 | U | 0.258 | 0.258 | r |
| 107-06-2 | 1,2-Dichloroethane(sim) | 0.247 | U | 0.247 | 0.247 | r |
| 71-55-6 | 1,1,1-Trichloroethane(sim) | 0.183 | U | 0.183 | 0.183 | r |
| 56-23-5 | Carbon Tetrachloride(sim) | 0.051 | J- | 0.032 | 0.032 | r |
| 75-35-4 | 1,1-Dichloroethene(sim) | 0.051 | U | 0.051 | 0.051 | r |
| 76-13-1 | Trichlorotrifluoroethane(sim) | 0.131 | U | 0.131 | 0.131 | r |
| 156-60-5 | Trans-1,2-Dichloroethene(sim) | 0.252 | UJ | 0.252 | 0.252 | r |
| 75-34-3 | 1,1-Dichloroethane(sim) | 0.247 | U | 0.247 | 0.247 | r |
| 156-59-2 | Cis-1,2-Dichloroethene(sim) | 0.051 | U | 0.051 | 0.051 | r |
| 67-66-3 | Chloroform(sim) | 0.205 | U | 0.205 | 0.205 | r |
| 78-87-5 | 1,2-dichloropropane(sim) | 0.217 | U | 0.217 | 0.217 | r |
| 75-27-4 | Bromodichloromethane(sim) | 0.149 | U | 0.149 | 0.149 | r |
| 79-01-6 | Trichloroethene(sim) | 0.037 | U | 0.037 | 0.037 | r |
| 123-91-1 | 1,4-Dioxane(sim) | 0.278 | U | 0.278 | 0.278 | r |
| 10061-01-5 | cis-1,3-Dichloropropene(sim) | 0.221 | U | 0.221 | 0.221 | r |
| 79-00-5 | 1,1,2-Trichloroethane(sim) | 0.183 | U | 0.183 | 0.183 | r |
| 124-48-1 | Dibromochloromethane(sim) | 0.118 | UJ | 0.118 | 0.118 | r |
| 106-93-4 | 1,2-Dibromoethane(EDB)(sim) | 0.130 | U | 0.130 | 0.130 | r |
| 127-18-4 | Tetrachloroethene(sim) | 0.145 | | 0.037 | 0.037 | r |
| 75-25-2 | Bromoform(sim) | 0.097 | UJ | 0.097 | 0.097 | r |
| 79-34-5 | 1,1,1,2-Tetrachloroethane(sim) | 0.146 | U | 0.146 | 0.146 | r |
| 100-44-7 | Benzyl chloride(sim) | 0.193 | U | 0.193 | 0.193 | r |
| 541-73-1 | 1,3-Dichlorobenzene(sim) | 0.166 | U | 0.166 | 0.166 | r |
| 106-46-7 | 1,4-Dichlorobenzene(sim) | 0.166 | U | 0.166 | 0.166 | r |
| 135-98-8 | sec-Butylbenzene(sim) | 0.182 | U | 0.182 | 0.182 | r |
| 99-87-6 | 4-Isopropyltoluene(sim) | 0.182 | U | 0.182 | 0.182 | r |
| 95-50-1 | 1,2-Dichlorobenzene(sim) | 0.166 | U | 0.166 | 0.166 | r |
| 104-51-8 | n-Butylbenzene(sim) | 0.182 | U | 0.182 | 0.182 | r |
| 120-82-1 | 1,2,4-Trichlorobenzene(sim) | 0.135 | U | 0.135 | 0.135 | r |
| 87-68-3 | Hexachlorobutadiene(sim) | 0.094 | U | 0.094 | 0.094 | r |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

JR 12/8/2025

1
AIR ANALYSIS DATA SHEET

CLIENT ID

OA-4

| | | | |
|----------------|----------|------------------|-------------------|
| Client: | ACT | Lab: | Phoenix Env. Labs |
| SDG No.: | GCP07983 | Lab Sample ID: | CP07984 |
| Canister: | 28588 | Lab File ID: | 0925_25.D |
| Instrument: | CHEM20 | Column: | RTX-1 60M |
| Date Received: | 09/25/23 | | |
| Purge Volume | 200 | (cc) | 09/25/23 |
| Date Analyzed: | 09/25/23 | | |
| Matrix: | AIR | Dilution Factor: | 1 |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|-------------|------------------------------------|-------|----|-------|-------|---|
| 115-07-1 | Propylene | 0.581 | U | 0.581 | 0.581 | r |
| 75-71-8 | Dichlorodifluoromethane | 0.503 | | 0.202 | 0.202 | r |
| 74-87-3 | Chloromethane | 0.526 | | 0.485 | 0.485 | r |
| 106-99-0 | 1,3-Butadiene | 0.452 | U | 0.452 | 0.452 | r |
| 75-00-3 | Chloroethane | 0.379 | U | 0.379 | 0.379 | r |
| 64-17-5 | Ethanol | 20.9 | J- | 0.531 | 0.531 | r |
| 67-64-1 | Acetone | 5.68 | J+ | 0.421 | 0.421 | r |
| 75-69-4 | Trichlorofluoromethane | 0.207 | | 0.178 | 0.178 | r |
| 67-63-0 | Isopropylalcohol | 0.841 | S | 0.407 | 0.407 | r |
| 107-13-1 | Acrylonitrile | 0.461 | U | 0.461 | 0.461 | r |
| 75-09-2 | Methylene Chloride | 0.863 | U | 0.863 | 0.863 | r |
| 75-15-0 | Carbon Disulfide | 0.321 | U | 0.321 | 0.321 | r |
| 1634-04-4 | Methyl tert-butyl ether(MTBE) | 0.278 | U | 0.278 | 0.278 | r |
| 78-93-3 | Methyl Ethyl Ketone | 0.339 | U | 0.339 | 0.339 | r |
| 110-54-3 | Hexane | 0.477 | S | 0.284 | 0.284 | r |
| 141-78-6 | Ethyl acetate | 0.278 | U | 0.278 | 0.278 | r |
| 109-99-9 | Tetrahydrofuran | 0.339 | U | 0.339 | 0.339 | r |
| 71-43-2 | Benzene | 0.329 | | 0.313 | 0.313 | r |
| 110-82-7 | Cyclohexane | 0.291 | U | 0.291 | 0.291 | r |
| 142-82-5 | Heptane | 0.244 | U | 0.244 | 0.244 | r |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | 0.244 | U | 0.244 | 0.244 | r |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.221 | U | 0.221 | 0.221 | r |
| 108-88-3 | Toluene | 0.833 | | 0.266 | 0.266 | r |
| 591-78-6 | 2-Hexanone(MBK) | 0.244 | U | 0.244 | 0.244 | r |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.146 | U | 0.146 | 0.146 | r |
| 108-90-7 | Chlorobenzene | 0.217 | U | 0.217 | 0.217 | r |
| 100-41-4 | Ethylbenzene | 0.230 | U | 0.230 | 0.230 | r |
| 179601-23-1 | m,p-Xylene | 0.494 | | 0.230 | 0.230 | r |
| 100-42-5 | Styrene | 0.235 | U | 0.235 | 0.235 | r |
| 95-47-6 | o-Xylene | 0.230 | U | 0.230 | 0.230 | r |
| 98-82-8 | Isopropylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 622-96-8 | 4-Ethyltoluene | 0.204 | U | 0.204 | 0.204 | r |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane(sim) | 0.143 | U | 0.143 | 0.143 | r |
| 75-01-4 | Vinyl Chloride(sim) | 0.078 | U | 0.078 | 0.078 | r |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

1
AIR ANALYSIS DATA SHEET

CLIENT ID

SS-18

| | | | |
|----------------|----------|------------------|-------------------|
| Client: | ACT | Lab: | Phoenix Env. Labs |
| SDG No.: | GCP07983 | Lab Sample ID: | CP07985 |
| Canister: | 28568 | Lab File ID: | 0925_28.D |
| Instrument: | CHEM20 | Column: | RTX-1 60M |
| Date Received: | 09/25/23 | | |
| Purge Volume | 200 (cc) | Date Analyzed: | 09/26/23 |
| Matrix: | AIR | Dilution Factor: | 1 |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|------------|------------------------------------|-------|----|-------|-------|---|
| 115-07-1 | Propylene | 0.581 | U | 0.581 | 0.581 | r |
| 75-71-8 | Dichlorodifluoromethane | 0.437 | | 0.202 | 0.202 | r |
| 74-87-3 | Chloromethane | 0.485 | U | 0.485 | 0.485 | r |
| 106-99-0 | 1,3-Butadiene | 0.452 | U | 0.452 | 0.452 | r |
| 75-00-3 | Chloroethane | 0.379 | U | 0.379 | 0.379 | r |
| 64-17-5 | Ethanol | 31.8 | J- | 0.531 | 0.531 | r |
| 67-64-1 | Acetone | 3.62 | J+ | 0.421 | 0.421 | r |
| 67-63-0 | Isopropylalcohol | 0.759 | S | 0.407 | 0.407 | r |
| 107-13-1 | Acrylonitrile | 0.461 | U | 0.461 | 0.461 | r |
| 75-09-2 | Methylene Chloride | 0.863 | U | 0.863 | 0.863 | r |
| 75-15-0 | Carbon Disulfide | 0.321 | U | 0.321 | 0.321 | r |
| 1634-04-4 | Methyl tert-butyl ether(MTBE) | 0.278 | U | 0.278 | 0.278 | r |
| 78-93-3 | Methyl Ethyl Ketone | 0.339 | U | 0.339 | 0.339 | r |
| 110-54-3 | Hexane | 0.284 | U | 0.284 | 0.284 | r |
| 141-78-6 | Ethyl acetate | 0.278 | U | 0.278 | 0.278 | r |
| 109-99-9 | Tetrahydrofuran | 0.339 | U | 0.339 | 0.339 | r |
| 110-82-7 | Cyclohexane | 0.291 | U | 0.291 | 0.291 | r |
| 142-82-5 | Heptane | 0.244 | U | 0.244 | 0.244 | r |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | 0.244 | U | 0.244 | 0.244 | r |
| 10061-02-6 | trans-1,3-Dichloropropene | 0.221 | U | 0.221 | 0.221 | r |
| 108-88-3 | Toluene | 0.506 | | 0.266 | 0.266 | r |
| 591-78-6 | 2-Hexanone(MBK) | 0.244 | U | 0.244 | 0.244 | r |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 0.146 | U | 0.146 | 0.146 | r |
| 108-90-7 | Chlorobenzene | 0.217 | U | 0.217 | 0.217 | r |
| 100-41-4 | Ethylbenzene | 0.230 | U | 0.230 | 0.230 | r |
| 100-42-5 | Styrene | 0.235 | U | 0.235 | 0.235 | r |
| 95-47-6 | o-Xylene | 0.230 | U | 0.230 | 0.230 | r |
| 98-82-8 | Isopropylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 622-96-8 | 4-Ethyltoluene | 0.204 | U | 0.204 | 0.204 | r |
| 108-67-8 | 1,3,5-Trimethylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 95-63-6 | 1,2,4-Trimethylbenzene | 0.204 | U | 0.204 | 0.204 | r |
| 106-46-7 | 1,4-Dichlorobenzene | 0.231 | | 0.166 | 0.166 | r |
| 76-14-2 | 1,2-Dichlorotetrafluoroethane(sim) | 0.143 | U | 0.143 | 0.143 | r |
| 75-01-4 | Vinyl Chloride(sim) | 0.078 | U | 0.078 | 0.078 | r |
| 74-83-9 | Bromomethane(sim) | 0.258 | U | 0.258 | 0.258 | r |
| 75-69-4 | Trichlorofluoromethane(sim) | 0.197 | | 0.178 | 0.178 | r |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

1
AIR ANALYSIS DATA SHEET

CLIENT ID

SS-18

| | | | |
|----------------|----------|------------------|-------------------|
| Client: | ACT | Lab: | Phoenix Env. Labs |
| SDG No.: | GCP07983 | Lab Sample ID: | CP07985 |
| Canister: | 28568 | Lab File ID: | 0925_28.D |
| Instrument: | CHEM20 | Column: | RTX-1 60M |
| Date Received: | 09/25/23 | | |
| Purge Volume | 200 | (cc) | 09/26/23 |
| Date Analyzed: | 09/26/23 | | |
| Matrix: | AIR | Dilution Factor: | 1 |

CONCENTRATION UNITS: (ppbv or ug/m3) ppbv

| CAS NO. | COMPOUND | CONC. | Q | MDL | PQL | R |
|-------------|--------------------------------|-------|----|-------|-------|---|
| 107-06-2 | 1,2-Dichloroethane(sim) | 0.247 | U | 0.247 | 0.247 | r |
| 71-55-6 | 1,1,1-Trichloroethane(sim) | 0.183 | U | 0.183 | 0.183 | r |
| 71-43-2 | Benzene(sim) | 0.313 | U | 0.313 | 0.313 | r |
| 56-23-5 | Carbon Tetrachloride(sim) | 0.040 | J- | 0.032 | 0.032 | r |
| 75-35-4 | 1,1-Dichloroethene(sim) | 0.051 | U | 0.051 | 0.051 | r |
| 76-13-1 | Trichlorotrifluoroethane(sim) | 0.131 | U | 0.131 | 0.131 | r |
| 156-60-5 | Trans-1,2-Dichloroethene(sim) | 0.252 | UJ | 0.252 | 0.252 | r |
| 75-34-3 | 1,1-Dichloroethane(sim) | 0.247 | U | 0.247 | 0.247 | r |
| 156-59-2 | Cis-1,2-Dichloroethene(sim) | 0.101 | | 0.051 | 0.051 | r |
| 67-66-3 | Chloroform(sim) | 0.205 | U | 0.205 | 0.205 | r |
| 78-87-5 | 1,2-dichloropropane(sim) | 0.217 | U | 0.217 | 0.217 | r |
| 75-27-4 | Bromodichloromethane(sim) | 0.149 | U | 0.149 | 0.149 | r |
| 79-01-6 | Trichloroethene(sim) | 0.177 | | 0.037 | 0.037 | r |
| 123-91-1 | 1,4-Dioxane(sim) | 0.278 | U | 0.278 | 0.278 | r |
| 10061-01-5 | cis-1,3-Dichloropropene(sim) | 0.221 | U | 0.221 | 0.221 | r |
| 79-00-5 | 1,1,2-Trichloroethane(sim) | 0.183 | U | 0.183 | 0.183 | r |
| 124-48-1 | Dibromochloromethane(sim) | 0.118 | UJ | 0.118 | 0.118 | r |
| 106-93-4 | 1,2-Dibromoethane(EDB)(sim) | 0.130 | U | 0.130 | 0.130 | r |
| 127-18-4 | Tetrachloroethene(sim) | 0.171 | | 0.037 | 0.037 | r |
| 75-25-2 | Bromoform(sim) | 0.097 | UJ | 0.097 | 0.097 | r |
| 179601-23-1 | m,p-Xylene(sim) | 0.371 | | 0.230 | 0.230 | r |
| 79-34-5 | 1,1,2,2-Tetrachloroethane(sim) | 0.146 | UJ | 0.146 | 0.146 | r |
| 100-44-7 | Benzyl chloride(sim) | 0.193 | U | 0.193 | 0.193 | r |
| 541-73-1 | 1,3-Dichlorobenzene(sim) | 0.166 | U | 0.166 | 0.166 | r |
| 135-98-8 | sec-Butylbenzene(sim) | 0.182 | U | 0.182 | 0.182 | r |
| 99-87-6 | 4-Isopropyltoluene(sim) | 0.182 | U | 0.182 | 0.182 | r |
| 95-50-1 | 1,2-Dichlorobenzene(sim) | 0.166 | U | 0.166 | 0.166 | r |
| 104-51-8 | n-Butylbenzene(sim) | 0.182 | U | 0.182 | 0.182 | r |
| 120-82-1 | 1,2,4-Trichlorobenzene(sim) | 0.135 | U | 0.135 | 0.135 | r |
| 87-68-3 | Hexachlorobutadiene(sim) | 0.094 | U | 0.094 | 0.094 | r |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |
| | | | | | | |

FORM I AIR

r=Result Reported U=Not Detected D=Reported Dilution E/J=Estimated Value X=Not Used S=Lab Solvent

JR 12/8/2025